



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:56 AM UTC

PDB ID : 3U3F / pdb_00003u3f
Title : Structural basis for the interaction of Pyk2 PAT domain with paxillin LD motifs
Authors : Vanarotti, M.; Miller, D.J.; Guibao, C.C.; Zheng, J.J.
Deposited on : 2011-10-05
Resolution : 3.10 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

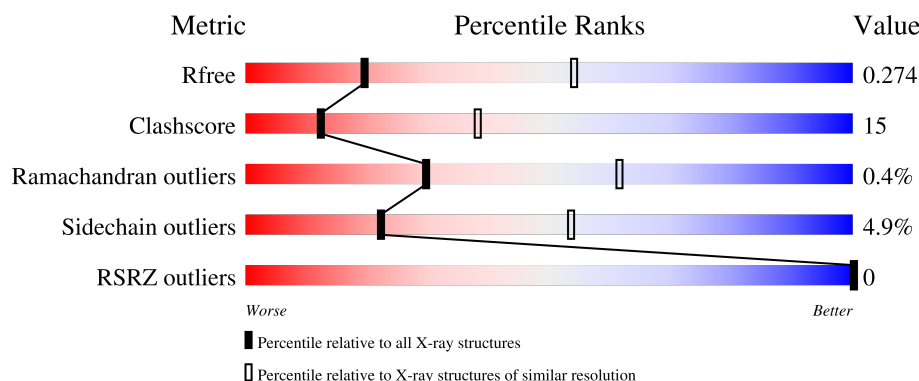
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

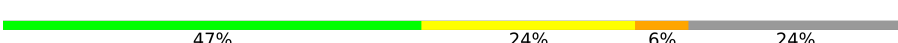
The reported resolution of this entry is 3.10 Å.

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.



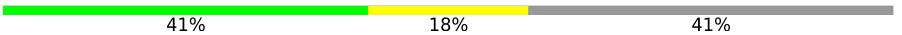
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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 $\leq 5\%$. 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	139	 60% 32% 6%
1	B	139	 59% 31% 7%
1	C	139	 57% 32% 9%
1	D	139	 60% 27% 9%
2	E	17	 47% 24% 6% 24%

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Mol	Chain	Length	Quality of chain
2	F	17	
2	G	17	
2	H	17	
2	I	17	
2	J	17	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4292 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 Protein-tyrosine kinase 2-beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	130	Total	C	N	O	S	0	0	0
			983	616	169	195	3			
1	B	129	Total	C	N	O	S	0	0	0
			978	615	166	194	3			
1	C	126	Total	C	N	O	S	0	0	0
			936	586	158	189	3			
1	D	126	Total	C	N	O	S	0	0	0
			937	588	158	188	3			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	867	GLY	-	expression tag	UNP Q14289
A	868	SER	-	expression tag	UNP Q14289
A	869	HIS	-	expression tag	UNP Q14289
A	870	MET	-	expression tag	UNP Q14289
A	899	SER	CYS	engineered mutation	UNP Q14289
B	867	GLY	-	expression tag	UNP Q14289
B	868	SER	-	expression tag	UNP Q14289
B	869	HIS	-	expression tag	UNP Q14289
B	870	MET	-	expression tag	UNP Q14289
B	899	SER	CYS	engineered mutation	UNP Q14289
C	867	GLY	-	expression tag	UNP Q14289
C	868	SER	-	expression tag	UNP Q14289
C	869	HIS	-	expression tag	UNP Q14289
C	870	MET	-	expression tag	UNP Q14289
C	899	SER	CYS	engineered mutation	UNP Q14289
D	867	GLY	-	expression tag	UNP Q14289
D	868	SER	-	expression tag	UNP Q14289
D	869	HIS	-	expression tag	UNP Q14289
D	870	MET	-	expression tag	UNP Q14289
D	899	SER	CYS	engineered mutation	UNP Q14289

- Molecule 2 is a protein called Paxillin LD2 peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	13	Total	C	N	O	S	0	0	0
			83	52	13	17	1			
2	F	13	Total	C	N	O	S	0	0	0
			83	52	13	17	1			
2	G	12	Total	C	N	O	S	0	0	0
			75	46	12	16	1			
2	H	14	Total	C	N	O	S	0	0	0
			88	55	14	18	1			
2	I	10	Total	C	N	O	S	0	0	0
			64	39	10	14	1			
2	J	10	Total	C	N	O	S	0	0	0
			62	37	10	14	1			

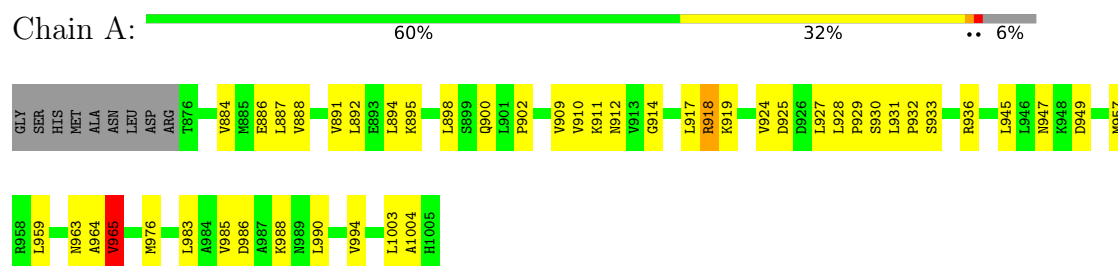
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	O	0	0
			2	2		
3	B	1	Total	O	0	0
			1	1		

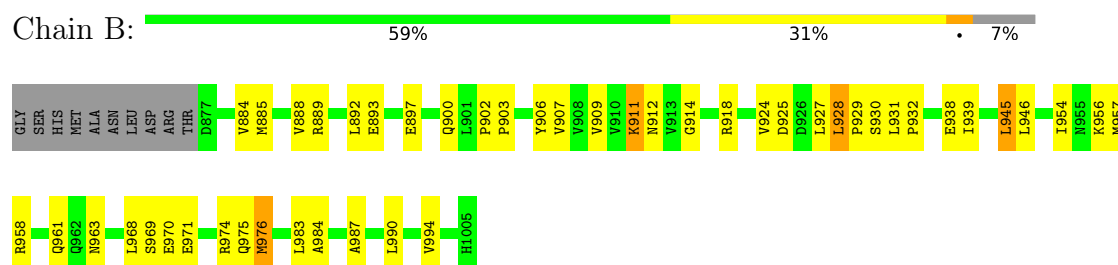
3 Residue-property plots [i](#)

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.

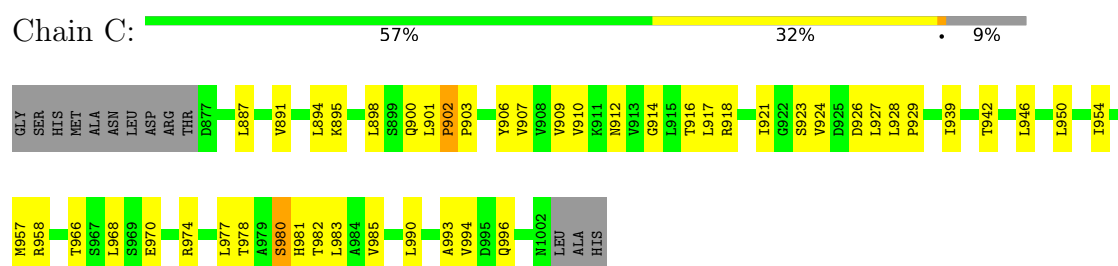
- Molecule 1: Protein-tyrosine kinase 2-beta



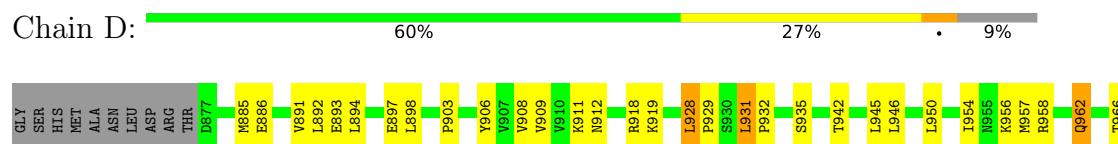
- Molecule 1: Protein-tyrosine kinase 2-beta

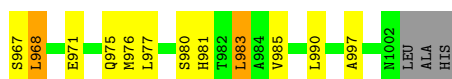


- Molecule 1: Protein-tyrosine kinase 2-beta



- Molecule 1: Protein-tyrosine kinase 2-beta





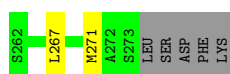
- Molecule 2: Paxillin LD2 peptide



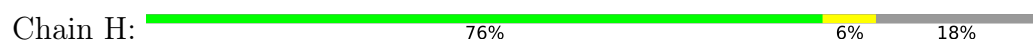
- Molecule 2: Paxillin LD2 peptide



- Molecule 2: Paxillin LD2 peptide



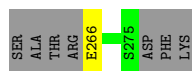
- Molecule 2: Paxillin LD2 peptide



- Molecule 2: Paxillin LD2 peptide



- Molecule 2: Paxillin LD2 peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.60Å 83.75Å 170.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.85 – 3.10 29.85 – 3.10	Depositor EDS
% Data completeness (in resolution range)	93.3 (29.85-3.10) 93.2 (29.85-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 3.11Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, R_{free}	0.218 , 0.266 0.217 , 0.274	Depositor DCC
R_{free} test set	1117 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	98.5	Xtriage
Anisotropy	0.098	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 115.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.469 for k,h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4292	wwPDB-VP
Average B, all atoms (Å ²)	108.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.14% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

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	0.52	0/989	0.89	3/1342 (0.2%)
1	B	0.55	0/984	0.91	2/1335 (0.1%)
1	C	0.49	0/942	0.87	3/1281 (0.2%)
1	D	0.54	0/943	0.91	1/1281 (0.1%)
2	E	0.65	0/82	0.97	0/111
2	F	0.56	0/82	0.94	0/111
2	G	0.48	0/74	0.99	0/100
2	H	0.49	0/87	0.87	0/118
2	I	0.35	0/63	0.82	0/84
2	J	0.37	0/61	0.80	0/81
All	All	0.52	0/4307	0.90	9/5844 (0.2%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	931	LEU	N-CA-C	6.89	114.33	108.13
1	B	902	PRO	CA-C-N	6.45	127.91	119.84
1	B	902	PRO	C-N-CA	6.45	127.91	119.84
1	A	902	PRO	CA-C-N	5.38	126.56	119.84
1	A	902	PRO	C-N-CA	5.38	126.56	119.84
1	C	907	VAL	CB-CA-C	-5.36	104.99	112.02
1	C	902	PRO	CA-C-N	5.26	126.41	119.84
1	C	902	PRO	C-N-CA	5.26	126.41	119.84
1	A	898	LEU	N-CA-C	-5.15	105.74	112.23

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	983	0	1019	34	0
1	B	978	0	1024	38	0
1	C	936	0	948	31	0
1	D	937	0	955	28	0
2	E	83	0	75	6	0
2	F	83	0	75	4	0
2	G	75	0	64	2	0
2	H	88	0	77	1	0
2	I	64	0	55	2	0
2	J	62	0	48	1	0
3	A	2	0	0	1	0
3	B	1	0	0	0	0
All	All	4292	0	4340	133	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 (133) 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:914:GLY:HA3	2:E:271:MET:HE1	1.43	0.97
1:A:931:LEU:HB2	1:A:932:PRO:HD2	1.54	0.88
1:B:970:GLU:H	1:B:970:GLU:CD	1.82	0.88
1:B:931:LEU:HB2	1:B:932:PRO:HD2	1.55	0.86
1:B:927:LEU:C	1:B:929:PRO:HD2	2.00	0.86
1:A:990:LEU:O	1:A:994:VAL:HG23	1.83	0.78
1:D:903:PRO:HA	1:D:906:TYR:CD1	2.19	0.77
1:C:891:VAL:HG13	1:C:980:SER:HB2	1.68	0.75
1:B:990:LEU:O	1:B:994:VAL:HG23	1.88	0.72
1:C:970:GLU:H	1:C:970:GLU:CD	1.99	0.71
1:D:886:GLU:OE2	1:D:919:LYS:HE3	1.91	0.70
1:D:891:VAL:HG13	1:D:980:SER:HB2	1.73	0.70
1:B:907:VAL:HG12	1:B:911:LYS:HE2	1.74	0.70
1:B:957:MET:HB2	1:B:976:MET:HE3	1.74	0.68
1:A:884:VAL:O	1:A:888:VAL:HG23	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:942:THR:OG1	1:C:990:LEU:HD12	1.95	0.66
1:C:903:PRO:HA	1:C:906:TYR:CD1	2.30	0.66
1:D:931:LEU:HB2	1:D:932:PRO:HD2	1.78	0.66
1:D:971:GLU:O	1:D:975:GLN:HG3	1.96	0.66
1:A:927:LEU:C	1:A:929:PRO:HD2	2.21	0.65
1:C:950:LEU:O	1:C:954:ILE:HG13	1.97	0.65
1:C:912:ASN:O	1:C:916:THR:HG23	1.97	0.64
1:A:933:SER:HA	1:A:936:ARG:HD3	1.79	0.64
1:C:993:ALA:O	1:C:996:GLN:HG2	1.98	0.64
1:B:884:VAL:O	1:B:888:VAL:HG23	1.98	0.63
1:A:886:GLU:HG2	3:A:3:HOH:O	1.99	0.62
1:A:925:ASP:HA	1:A:928:LEU:HD23	1.82	0.62
1:D:903:PRO:HA	1:D:906:TYR:CE1	2.34	0.62
1:B:970:GLU:CD	1:B:970:GLU:N	2.53	0.61
1:B:945:LEU:HD12	1:B:945:LEU:O	2.02	0.60
1:B:928:LEU:N	1:B:929:PRO:HD2	2.16	0.60
1:C:978:THR:O	1:C:982:THR:HG23	2.02	0.59
1:B:914:GLY:HA3	2:F:271:MET:HE1	1.85	0.59
1:A:928:LEU:N	1:A:929:PRO:HD2	2.18	0.59
2:F:271:MET:HE3	2:F:274:LEU:HD12	1.83	0.59
1:C:887:LEU:HD12	1:C:917:LEU:HA	1.86	0.57
1:C:898:LEU:HD12	1:C:977:LEU:HD21	1.86	0.57
1:B:918:ARG:HG3	2:F:267:LEU:HD23	1.86	0.57
1:A:909:VAL:O	1:A:912:ASN:HB2	2.04	0.57
1:A:930:SER:O	1:A:931:LEU:HB3	2.05	0.56
1:D:928:LEU:O	1:D:931:LEU:HG	2.06	0.56
1:C:903:PRO:HA	1:C:906:TYR:CE1	2.41	0.55
1:D:983:LEU:HD12	1:D:983:LEU:O	2.07	0.54
1:A:1003:LEU:HD12	1:A:1004:ALA:N	2.22	0.54
1:D:928:LEU:N	1:D:929:PRO:HD2	2.23	0.53
1:A:964:ALA:C	1:A:965:VAL:HG23	2.33	0.53
1:D:950:LEU:O	1:D:954:ILE:HG13	2.09	0.53
1:D:966:THR:C	1:D:968:LEU:H	2.17	0.53
1:B:930:SER:O	1:B:931:LEU:HB3	2.08	0.53
1:B:924:VAL:HG21	1:B:990:LEU:HD21	1.89	0.52
1:B:925:ASP:HA	1:B:928:LEU:HD23	1.90	0.52
1:A:947:ASN:OD1	2:E:267:LEU:HD13	2.09	0.52
1:D:935:SER:HB2	1:D:997:ALA:HB2	1.91	0.51
1:A:924:VAL:HG21	1:A:990:LEU:HD21	1.92	0.51
1:A:957:MET:HB2	1:A:976:MET:HE3	1.91	0.51
1:C:894:LEU:C	1:C:894:LEU:HD23	2.36	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:897:GLU:HA	1:B:900:GLN:HG2	1.92	0.51
1:A:909:VAL:HG23	1:A:910:VAL:N	2.26	0.50
1:B:909:VAL:O	1:B:912:ASN:HB2	2.11	0.50
1:A:917:LEU:HD12	1:A:917:LEU:O	2.12	0.49
1:D:898:LEU:HD12	1:D:977:LEU:HD21	1.94	0.49
1:A:919:LYS:HD2	1:C:918:ARG:NE	2.28	0.49
1:B:928:LEU:N	1:B:929:PRO:CD	2.76	0.49
1:D:908:VAL:O	1:D:912:ASN:ND2	2.43	0.48
2:E:271:MET:O	2:E:274:LEU:HB2	2.15	0.47
1:C:939:ILE:HD13	1:C:994:VAL:HG22	1.96	0.47
1:D:958:ARG:O	1:D:962:GLN:HG2	2.14	0.47
1:D:981:HIS:O	1:D:985:VAL:HG23	2.14	0.47
1:B:938:GLU:HA	1:B:938:GLU:OE1	2.15	0.46
1:D:894:LEU:C	1:D:894:LEU:HD23	2.40	0.46
1:A:959:LEU:O	1:A:963:ASN:HB2	2.16	0.46
1:B:903:PRO:HA	1:B:906:TYR:CE1	2.50	0.46
1:D:893:GLU:HB3	1:D:897:GLU:OE1	2.16	0.46
1:B:931:LEU:HB2	1:B:932:PRO:CD	2.38	0.46
1:D:892:LEU:HA	1:D:892:LEU:HD23	1.72	0.45
1:A:918:ARG:HG3	2:E:267:LEU:HD23	1.98	0.45
1:A:985:VAL:O	1:A:988:LYS:HB3	2.16	0.45
1:C:887:LEU:C	1:C:887:LEU:HD23	2.42	0.45
1:D:897:GLU:OE2	1:D:909:VAL:HG13	2.17	0.45
1:C:901:LEU:HD22	1:C:902:PRO:HD2	1.98	0.45
1:B:971:GLU:O	1:B:975:GLN:HG3	2.17	0.44
1:C:946:LEU:CD2	1:C:983:LEU:HD12	2.48	0.44
2:I:267:LEU:O	2:I:271:MET:HB2	2.16	0.44
1:B:885:MET:HB3	1:B:885:MET:HE3	1.76	0.44
1:C:921:ILE:HD12	2:G:267:LEU:HD22	1.99	0.44
1:C:981:HIS:O	1:C:985:VAL:HG23	2.18	0.44
1:B:954:ILE:HD12	2:F:270:LEU:HD22	1.98	0.44
1:D:962:GLN:HG2	1:D:962:GLN:H	1.55	0.44
1:A:1003:LEU:HD12	1:A:1003:LEU:C	2.43	0.44
1:C:966:THR:C	1:C:968:LEU:H	2.26	0.44
1:B:974:ARG:HH11	1:B:974:ARG:HG2	1.83	0.44
1:A:892:LEU:HD21	2:I:274:LEU:HA	1.99	0.44
1:C:928:LEU:N	1:C:929:PRO:HD2	2.33	0.43
1:A:911:LYS:HE3	2:E:274:LEU:HB3	2.00	0.43
1:C:914:GLY:HA3	2:G:271:MET:HE1	2.00	0.43
1:C:910:VAL:HG21	1:C:957:MET:HG3	2.00	0.43
1:A:945:LEU:HG	1:A:986:ASP:OD2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:984:ALA:O	1:B:987:ALA:HB3	2.18	0.43
1:B:924:VAL:O	1:B:928:LEU:HD22	2.19	0.43
1:D:946:LEU:HD22	1:D:983:LEU:CD1	2.49	0.43
1:B:889:ARG:O	1:B:893:GLU:HG3	2.19	0.43
1:B:903:PRO:HA	1:B:906:TYR:CD1	2.54	0.43
1:A:894:LEU:C	1:A:894:LEU:HD23	2.44	0.42
1:A:945:LEU:HD12	1:A:945:LEU:O	2.19	0.42
1:A:919:LYS:HD2	1:C:918:ARG:HE	1.84	0.42
1:C:894:LEU:HD12	1:C:909:VAL:HG12	2.01	0.42
1:D:942:THR:O	1:D:945:LEU:HB3	2.20	0.42
1:B:963:ASN:OD1	1:B:968:LEU:HD11	2.19	0.42
1:C:927:LEU:O	1:C:928:LEU:C	2.61	0.42
1:D:966:THR:O	1:D:968:LEU:N	2.48	0.42
1:B:931:LEU:C	1:B:931:LEU:HD12	2.44	0.42
1:B:969:SER:HB3	1:B:970:GLU:OE2	2.20	0.42
1:C:970:GLU:O	1:C:974:ARG:HB2	2.20	0.42
1:D:942:THR:OG1	1:D:990:LEU:HD12	2.19	0.42
1:B:946:LEU:HD22	1:B:983:LEU:HD12	2.02	0.41
1:C:895:LYS:HA	1:C:977:LEU:CD2	2.50	0.41
1:C:923:SER:O	1:C:926:ASP:HB2	2.20	0.41
1:A:891:VAL:O	1:A:895:LYS:HG3	2.19	0.41
1:D:918:ARG:HB2	2:H:267:LEU:HD21	2.02	0.41
1:B:945:LEU:HD12	1:B:945:LEU:C	2.45	0.41
1:C:917:LEU:HD23	1:C:950:LEU:HD22	2.01	0.41
1:D:957:MET:HB2	1:D:976:MET:HE3	2.02	0.41
1:B:907:VAL:CG2	1:B:961:GLN:HE22	2.33	0.41
1:A:949:ASP:HB3	1:A:983:LEU:HD13	2.03	0.41
1:C:895:LYS:HA	1:C:977:LEU:HD22	2.03	0.41
1:A:887:LEU:HD12	1:A:917:LEU:HA	2.02	0.41
1:A:892:LEU:HD23	1:A:892:LEU:HA	1.85	0.41
1:B:892:LEU:HA	1:B:892:LEU:HD23	1.82	0.41
1:B:927:LEU:HD12	1:B:927:LEU:HA	1.87	0.41
2:J:266:GLU:HA	2:J:266:GLU:OE2	2.21	0.41
2:E:264:THR:O	2:E:267:LEU:HB3	2.21	0.40
1:A:909:VAL:CG2	1:A:910:VAL:N	2.84	0.40
1:B:893:GLU:OE1	1:D:911:LYS:NZ	2.55	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	128/139 (92%)	119 (93%)	8 (6%)	1 (1%)	16	47
1	B	127/139 (91%)	121 (95%)	6 (5%)	0	100	100
1	C	124/139 (89%)	118 (95%)	6 (5%)	0	100	100
1	D	124/139 (89%)	117 (94%)	6 (5%)	1 (1%)	16	47
2	E	11/17 (65%)	10 (91%)	1 (9%)	0	100	100
2	F	11/17 (65%)	10 (91%)	1 (9%)	0	100	100
2	G	10/17 (59%)	8 (80%)	2 (20%)	0	100	100
2	H	12/17 (71%)	10 (83%)	2 (17%)	0	100	100
2	I	8/17 (47%)	8 (100%)	0	0	100	100
2	J	8/17 (47%)	8 (100%)	0	0	100	100
All	All	563/658 (86%)	529 (94%)	32 (6%)	2 (0%)	30	61

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	967	SER
1	A	965	VAL

5.3.2 Protein sidechains

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	110/122 (90%)	107 (97%)	3 (3%)	39	67

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	111/122 (91%)	104 (94%)	7 (6%)	16	45
1	C	103/122 (84%)	99 (96%)	4 (4%)	28	60
1	D	103/122 (84%)	97 (94%)	6 (6%)	18	48
2	E	7/15 (47%)	5 (71%)	2 (29%)	0	1
2	F	7/15 (47%)	6 (86%)	1 (14%)	3	14
2	G	6/15 (40%)	6 (100%)	0	100	100
2	H	7/15 (47%)	7 (100%)	0	100	100
2	I	6/15 (40%)	6 (100%)	0	100	100
2	J	5/15 (33%)	5 (100%)	0	100	100
All	All	465/578 (80%)	442 (95%)	23 (5%)	22	53

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	900	GLN
1	A	918	ARG
1	A	965	VAL
1	B	911	LYS
1	B	928	LEU
1	B	939	ILE
1	B	945	LEU
1	B	956	LYS
1	B	958	ARG
1	B	976	MET
1	C	900	GLN
1	C	924	VAL
1	C	958	ARG
1	C	980	SER
1	D	885	MET
1	D	928	LEU
1	D	956	LYS
1	D	962	GLN
1	D	968	LEU
1	D	983	LEU
2	E	270	LEU
2	E	274	LEU
2	F	271	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	883	ASN
1	B	955	ASN
1	B	996	GLN
1	C	883	ASN
1	C	963	ASN
1	D	883	ASN
1	D	912	ASN
1	D	961	GLN
1	D	981	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 oligosaccharides 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	130/139 (93%)	-1.34	0 100 100	66, 99, 137, 163	0
1	B	129/139 (92%)	-1.26	0 100 100	68, 99, 136, 157	0
1	C	126/139 (90%)	-1.25	0 100 100	72, 102, 157, 180	0
1	D	126/139 (90%)	-1.26	0 100 100	73, 104, 154, 183	0
2	E	13/17 (76%)	-1.03	0 100 100	95, 120, 142, 144	0
2	F	13/17 (76%)	-1.21	0 100 100	95, 120, 152, 154	0
2	G	12/17 (70%)	-1.19	0 100 100	97, 112, 144, 166	0
2	H	14/17 (82%)	-1.23	0 100 100	98, 114, 154, 165	0
2	I	10/17 (58%)	-0.88	0 100 100	126, 146, 152, 167	0
2	J	10/17 (58%)	-1.05	0 100 100	139, 151, 156, 158	0
All	All	583/658 (88%)	-1.26	0 100 100	66, 104, 154, 183	0

There are no RSRZ outliers to report.

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 oligosaccharides 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.