



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:52 AM UTC

PDB ID : 1OW6 / pdb_00001ow6
Title : Paxillin LD4 motif bound to the Focal Adhesion Targeting (FAT) domain of the Focal Adhesion Kinase
Authors : Hoellerer, M.K.; Noble, M.E.M.; Labesse, G.; Campbell, I.D.; Werner, J.M.; Arold, S.T.
Deposited on : 2003-03-28
Resolution : 2.35 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

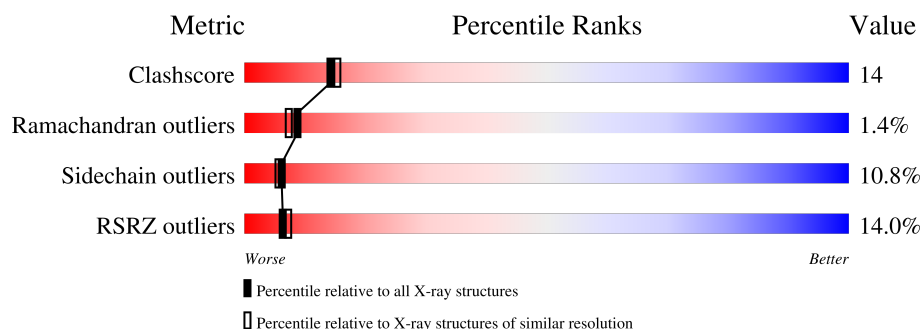
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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))
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	161	
1	B	161	
1	C	161	
2	D	13	
2	F	13	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3488 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 Focal adhesion kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	134	Total	C	N	O	S	0	0	0
			1042	658	176	201	7			
1	B	139	Total	C	N	O	S	0	0	0
			1073	680	180	206	7			
1	C	139	Total	C	N	O	S	0	0	0
			1073	680	180	206	7			

- Molecule 2 is a protein called Paxillin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	12	Total	C	N	O	S	0	0	0
			94	56	15	22	1			
2	F	11	Total	C	N	O	S	0	0	0
			87	53	14	19	1			

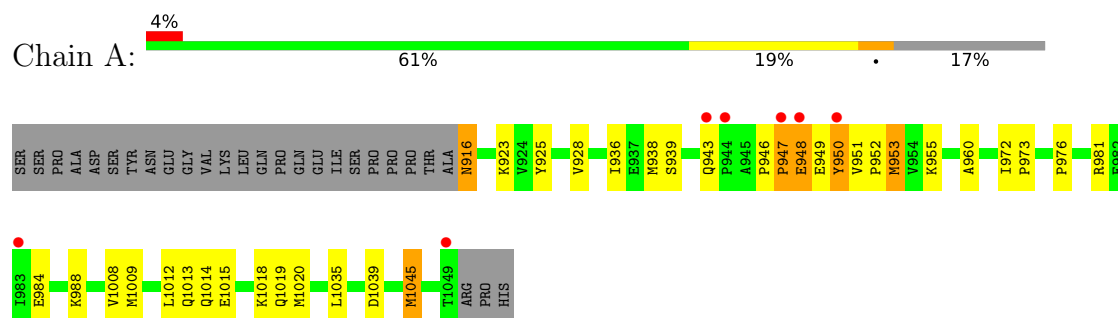
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	36	Total	O	0	0
			36	36		
3	B	44	Total	O	0	0
			44	44		
3	C	35	Total	O	0	0
			35	35		
3	D	3	Total	O	0	0
			3	3		
3	F	1	Total	O	0	0
			1	1		

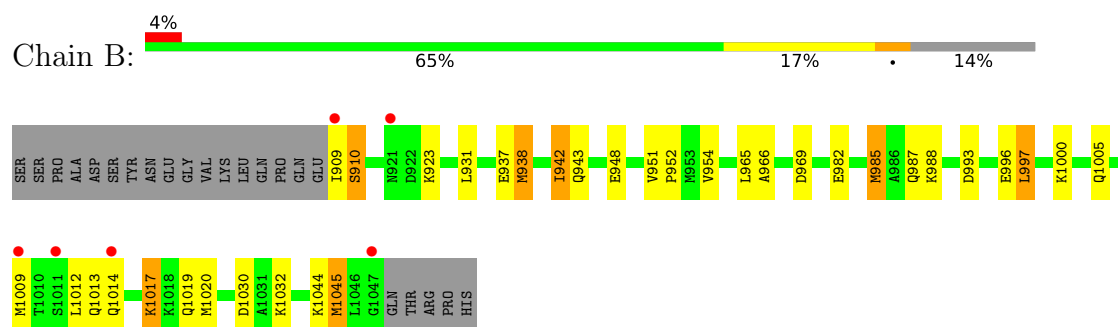
3 Residue-property plots [i](#)

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.

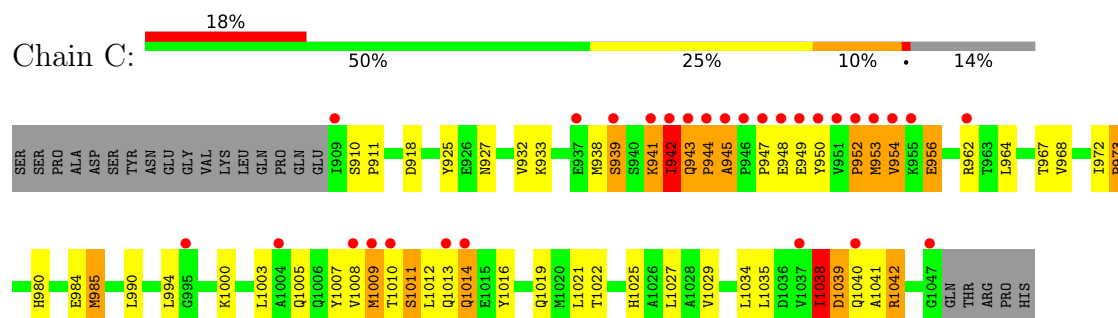
- Molecule 1: Focal adhesion kinase 1



- Molecule 1: Focal adhesion kinase 1



- Molecule 1: Focal adhesion kinase 1

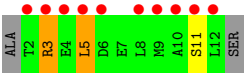
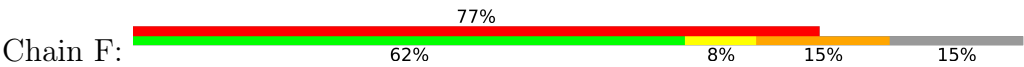


- Molecule 2: Paxillin





● Molecule 2: Paxillin



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	88.31Å 223.27Å 96.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.27 – 2.35 37.27 – 2.35	Depositor EDS
% Data completeness (in resolution range)	96.8 (37.27-2.35) 96.8 (37.27-2.35)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.242 , 0.274 0.244 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	44.2	Xtriage
Anisotropy	0.722	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 62.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3488	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.38% 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	0.68	0/1053	1.39	4/1424 (0.3%)
1	B	0.72	0/1087	1.41	6/1474 (0.4%)
1	C	0.76	2/1087 (0.2%)	1.56	11/1474 (0.7%)
2	D	0.50	0/93	1.47	1/122 (0.8%)
2	F	0.78	0/86	1.34	1/114 (0.9%)
All	All	0.72	2/3406 (0.1%)	1.45	23/4608 (0.5%)

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	C	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	1038	ILE	CA-CB	10.75	1.67	1.54
1	C	1038	ILE	CG1-CD1	5.24	1.72	1.51

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	1038	ILE	N-CA-C	9.44	121.38	110.62
1	C	1040	GLN	N-CA-C	-9.43	100.87	112.38
1	C	1042	ARG	N-CA-C	-9.26	102.49	113.88
1	C	1039	ASP	N-CA-C	-8.63	101.95	111.36
1	B	931	LEU	N-CA-C	-7.00	104.36	113.12
1	B	1005	GLN	N-CA-C	-6.90	103.69	111.14
1	A	928	VAL	CB-CA-C	-6.43	103.34	112.22
1	C	984	GLU	N-CA-C	-6.06	104.75	111.36
1	A	1039	ASP	N-CA-C	-5.84	105.00	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	960	ALA	N-CA-C	-5.80	105.09	111.82
1	B	937	GLU	N-CA-C	-5.79	104.89	111.14
1	C	932	VAL	CB-CA-C	-5.76	104.60	111.97
1	C	942	ILE	N-CA-C	5.76	121.31	109.34
1	B	997	LEU	N-CA-C	-5.58	105.20	112.23
2	D	8	LEU	N-CA-C	5.56	117.15	111.14
1	A	928	VAL	N-CA-C	5.51	116.28	110.72
1	B	942	ILE	N-CA-C	5.45	119.67	112.04
1	C	925	TYR	N-CA-C	-5.39	105.30	111.07
2	F	3	ARG	N-CA-C	5.36	119.59	112.30
1	C	911	PRO	N-CA-C	-5.34	104.19	110.70
1	C	945	ALA	N-CA-C	5.32	121.58	109.81
1	B	1017	LYS	N-CA-C	-5.12	105.78	111.36
1	C	939	SER	N-CA-C	5.09	118.27	111.75

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	1038	ILE	Mainchain

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	1042	0	1093	32	0
1	B	1073	0	1127	23	0
1	C	1073	0	1127	42	0
2	D	94	0	92	3	0
2	F	87	0	87	1	0
3	A	36	0	0	2	0
3	B	44	0	0	3	0
3	C	35	0	0	0	0
3	D	3	0	0	1	0
3	F	1	0	0	0	0
All	All	3488	0	3526	98	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 14.

All (98) 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:947:PRO:HA	1:A:950:TYR:CE2	2.12	0.84
1:A:947:PRO:HA	1:A:950:TYR:CZ	2.16	0.81
1:C:1007:TYR:O	1:C:1013:GLN:HB2	1.83	0.78
1:A:1009:MET:H	1:A:1013:GLN:NE2	1.81	0.77
1:A:976:PRO:HD3	1:A:1045:MET:HG2	1.67	0.76
1:B:965:LEU:HD22	1:B:987:GLN:HG2	1.68	0.76
1:B:1045:MET:HE3	1:B:1045:MET:HA	1.68	0.75
1:B:938:MET:HE1	1:B:997:LEU:HD11	1.69	0.75
1:B:985:MET:HE3	1:B:988:LYS:HD2	1.68	0.74
1:B:1013:GLN:HB3	3:B:40:HOH:O	1.89	0.72
1:A:1009:MET:H	1:A:1013:GLN:HE21	1.37	0.72
1:B:985:MET:CE	1:B:988:LYS:HD2	2.20	0.70
1:C:943:GLN:HB3	1:C:944:PRO:HD3	1.74	0.70
1:C:943:GLN:HB3	1:C:944:PRO:CD	2.24	0.67
1:C:1010:THR:HG22	1:C:1012:LEU:H	1.60	0.67
1:B:951:VAL:HB	1:B:952:PRO:HD3	1.79	0.64
1:B:993:ASP:OD2	1:B:1030:ASP:OD2	2.21	0.59
1:A:951:VAL:HG12	1:A:955:LYS:HE3	1.84	0.59
1:B:938:MET:HE2	1:B:954:VAL:HG22	1.84	0.59
1:C:1038:ILE:O	1:C:1042:ARG:HG3	2.02	0.58
1:A:1045:MET:HA	1:A:1045:MET:HE2	1.83	0.58
1:A:947:PRO:HB3	1:A:950:TYR:OH	2.03	0.58
1:B:965:LEU:CD2	1:B:987:GLN:HG2	2.34	0.56
1:A:946:PRO:HG2	1:A:949:GLU:HB2	1.87	0.56
1:C:1025:HIS:O	1:C:1029:VAL:HG23	2.05	0.56
1:B:1012:LEU:HG	1:C:985:MET:HG3	1.88	0.56
1:A:1014:GLN:HE21	1:A:1018:LYS:NZ	2.04	0.55
1:A:949:GLU:HA	3:A:93:HOH:O	2.05	0.55
1:A:950:TYR:CB	1:A:953:MET:HE2	2.37	0.55
1:C:943:GLN:CB	1:C:944:PRO:CD	2.85	0.54
1:C:942:ILE:HG13	1:C:950:TYR:HB3	1.90	0.54
2:D:4:GLU:HG3	3:D:37:HOH:O	2.06	0.54
1:B:909:ILE:HD12	3:B:19:HOH:O	2.08	0.53
1:B:1000:LYS:HD2	1:B:1019:GLN:HB3	1.90	0.53
1:A:950:TYR:HB3	1:A:953:MET:HE2	1.91	0.53
1:C:945:ALA:HB1	1:C:949:GLU:HB2	1.93	0.51
1:A:1045:MET:HA	1:A:1045:MET:CE	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:951:VAL:HB	1:A:952:PRO:HD3	1.92	0.50
1:C:927:ASN:ND2	1:C:967:THR:OG1	2.45	0.50
1:C:1039:ASP:C	1:C:1041:ALA:N	2.63	0.50
1:A:1014:GLN:HE21	1:A:1018:LYS:CE	2.25	0.49
1:C:941:LYS:O	1:C:943:GLN:N	2.45	0.49
1:A:1008:VAL:HG22	1:A:1009:MET:HG2	1.94	0.48
1:C:968:VAL:HG21	1:C:1034:LEU:HD21	1.95	0.48
1:A:946:PRO:O	1:A:947:PRO:C	2.57	0.47
1:A:951:VAL:N	1:A:952:PRO:CD	2.77	0.47
1:A:972:ILE:N	1:A:973:PRO:CD	2.76	0.47
1:A:976:PRO:HD3	1:A:1045:MET:CG	2.42	0.47
1:C:1003:LEU:HD13	1:C:1016:TYR:CE1	2.49	0.47
1:A:916:ASN:OD1	1:A:916:ASN:N	2.48	0.47
1:C:1014:GLN:O	1:C:1014:GLN:NE2	2.47	0.47
1:A:943:GLN:O	3:A:73:HOH:O	2.20	0.47
1:B:985:MET:HE1	1:B:988:LYS:HE3	1.96	0.47
1:C:918:ASP:OD1	1:C:918:ASP:C	2.58	0.47
1:C:985:MET:HA	1:C:985:MET:HE3	1.96	0.47
1:B:1014:GLN:HG2	3:B:59:HOH:O	2.13	0.46
1:C:990:LEU:HD23	1:C:1027:LEU:HD12	1.98	0.46
1:B:966:ALA:O	1:B:969:ASP:HB2	2.16	0.46
1:C:1038:ILE:C	1:C:1041:ALA:H	2.23	0.46
1:A:939:SER:O	1:A:943:GLN:HB2	2.15	0.45
1:C:947:PRO:HB3	1:C:1005:GLN:HE21	1.82	0.45
1:C:1038:ILE:O	1:C:1038:ILE:HG22	2.15	0.45
1:C:962:ARG:HG3	2:F:5:LEU:HD21	1.98	0.45
1:B:985:MET:HG3	1:C:1012:LEU:HG	1.98	0.45
1:B:938:MET:HE2	1:B:954:VAL:CG2	2.47	0.45
1:C:1009:MET:HE2	1:C:1009:MET:HB2	1.91	0.44
2:D:5:LEU:HA	2:D:8:LEU:HD12	1.99	0.44
1:A:947:PRO:CA	1:A:950:TYR:CZ	2.94	0.44
1:C:1010:THR:HG22	1:C:1011:SER:N	2.33	0.44
1:C:972:ILE:HG22	1:C:980:HIS:CE1	2.53	0.44
2:D:3:ARG:O	2:D:7:GLU:HG2	2.18	0.43
1:C:1012:LEU:HD13	1:C:1016:TYR:HE2	1.82	0.43
1:C:972:ILE:HB	1:C:973:PRO:HD3	2.00	0.43
1:B:996:GLU:O	1:B:1000:LYS:HG2	2.19	0.43
1:B:909:ILE:HG13	1:B:910:SER:N	2.34	0.43
1:C:985:MET:HA	1:C:985:MET:CE	2.49	0.42
1:C:1039:ASP:C	1:C:1041:ALA:H	2.27	0.42
1:C:939:SER:HA	1:C:1021:LEU:HD22	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:938:MET:O	1:B:942:ILE:HB	2.19	0.42
1:C:972:ILE:N	1:C:973:PRO:CD	2.83	0.42
1:A:946:PRO:O	1:A:948:GLU:N	2.52	0.42
1:B:985:MET:HE3	1:B:985:MET:HA	2.02	0.42
1:C:1000:LYS:HG3	1:C:1019:GLN:HB3	2.02	0.41
1:C:952:PRO:C	1:C:954:VAL:N	2.75	0.41
1:C:990:LEU:CD2	1:C:1027:LEU:HD12	2.50	0.41
1:C:954:VAL:C	1:C:956:GLU:N	2.78	0.41
1:A:950:TYR:HB3	1:A:953:MET:CE	2.50	0.41
1:B:1017:LYS:O	1:B:1020:MET:HB2	2.21	0.41
1:A:1009:MET:N	1:A:1013:GLN:HE21	2.13	0.41
1:A:925:TYR:CD1	1:A:925:TYR:C	2.99	0.41
1:A:1008:VAL:HA	1:A:1013:GLN:HE21	1.85	0.41
1:C:942:ILE:O	1:C:943:GLN:C	2.64	0.41
1:C:942:ILE:HG23	1:C:943:GLN:N	2.35	0.41
1:C:964:LEU:O	1:C:968:VAL:HG23	2.21	0.41
1:A:1015:GLU:O	1:A:1019:GLN:HG3	2.21	0.41
1:C:938:MET:HA	1:C:953:MET:SD	2.61	0.40
1:C:994:LEU:HD12	1:C:994:LEU:HA	1.96	0.40
1:A:936:ILE:HD13	1:A:936:ILE:HG21	1.79	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	132/161 (82%)	127 (96%)	4 (3%)	1 (1%)	16	17
1	B	137/161 (85%)	136 (99%)	1 (1%)	0	100	100
1	C	137/161 (85%)	121 (88%)	11 (8%)	5 (4%)	2	1
2	D	10/13 (77%)	10 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	F	9/13 (69%)	8 (89%)	1 (11%)	0	100	100
All	All	425/509 (84%)	402 (95%)	17 (4%)	6 (1%)	9	7

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	942	ILE
1	C	943	GLN
1	C	953	MET
1	A	947	PRO
1	C	944	PRO
1	C	952	PRO

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	117/141 (83%)	104 (89%)	13 (11%)	6	5
1	B	121/141 (86%)	110 (91%)	11 (9%)	9	9
1	C	121/141 (86%)	107 (88%)	14 (12%)	5	5
2	D	11/11 (100%)	11 (100%)	0	100	100
2	F	10/11 (91%)	7 (70%)	3 (30%)	0	0
All	All	380/445 (85%)	339 (89%)	41 (11%)	6	5

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	916	ASN
1	A	923	LYS
1	A	938	MET
1	A	948	GLU
1	A	950	TYR
1	A	953	MET

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Mol	Chain	Res	Type
1	A	981	ARG
1	A	984	GLU
1	A	988	LYS
1	A	1012	LEU
1	A	1020	MET
1	A	1035	LEU
1	A	1045	MET
1	B	910	SER
1	B	923	LYS
1	B	938	MET
1	B	943	GLN
1	B	948	GLU
1	B	982	GLU
1	B	985	MET
1	B	1009	MET
1	B	1032	LYS
1	B	1044	LYS
1	B	1045	MET
1	C	910	SER
1	C	933	LYS
1	C	941	LYS
1	C	948	GLU
1	C	954	VAL
1	C	956	GLU
1	C	973	PRO
1	C	985	MET
1	C	1008	VAL
1	C	1009	MET
1	C	1011	SER
1	C	1014	GLN
1	C	1022	THR
1	C	1035	LEU
2	F	3	ARG
2	F	5	LEU
2	F	11	SER

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	916	ASN
1	A	921	ASN
1	A	1013	GLN

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Mol	Chain	Res	Type
1	A	1014	GLN
1	B	921	ASN
1	C	927	ASN
1	C	1005	GLN
1	C	1014	GLN
1	C	1033	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 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

6.1 Protein, DNA and RNA chains

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	134/161 (83%)	0.83	7 (5%) 33 38	37, 54, 73, 87	0
1	B	139/161 (86%)	0.49	6 (4%) 40 45	30, 48, 63, 70	0
1	C	139/161 (86%)	1.32	29 (20%) 2 2	20, 56, 108, 116	0
2	D	12/13 (92%)	2.63	9 (75%) 0 0	70, 89, 103, 105	0
2	F	11/13 (84%)	5.59	10 (90%) 0 0	131, 132, 135, 137	0
All	All	435/509 (85%)	1.05	61 (14%) 6 7	20, 53, 106, 137	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	2	THR	9.0
2	F	12	LEU	8.9
2	F	9	MET	8.3
2	F	5	LEU	8.0
2	F	8	LEU	7.7
1	C	942	ILE	7.0
1	C	951	VAL	6.4
2	F	10	ALA	6.2
1	C	1009	MET	6.1
1	C	945	ALA	4.8
1	C	953	MET	4.4
1	C	950	TYR	4.4
1	C	1047	GLY	4.3
1	C	944	PRO	4.2
1	A	950	TYR	4.2
2	D	5	LEU	4.0
1	C	943	GLN	4.0
1	A	943	GLN	3.8
2	D	2	THR	3.8
2	D	12	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
1	C	1004	ALA	3.7
2	F	6	ASP	3.6
2	D	13	SER	3.5
1	C	941	LYS	3.5
2	F	11	SER	3.3
1	A	947	PRO	3.3
1	A	1049	THR	3.3
1	C	954	VAL	3.1
1	C	946	PRO	2.9
1	C	947	PRO	2.8
1	C	909	ILE	2.7
1	C	1013	GLN	2.7
2	D	8	LEU	2.7
2	D	10	ALA	2.7
1	C	955	LYS	2.6
1	A	948	GLU	2.6
1	C	1010	THR	2.6
1	B	921	ASN	2.5
2	D	3	ARG	2.5
1	B	1009	MET	2.4
2	F	4	GLU	2.4
1	C	952	PRO	2.4
1	C	948	GLU	2.4
1	C	995	GLY	2.3
2	D	6	ASP	2.3
1	B	1011	SER	2.3
2	F	3	ARG	2.3
1	B	1047	GLY	2.3
1	C	937	GLU	2.3
1	C	1008	VAL	2.3
2	D	11	SER	2.2
1	B	909	ILE	2.2
1	B	1014	GLN	2.2
1	A	983	ILE	2.2
1	C	949	GLU	2.2
1	C	962	ARG	2.2
1	C	1037	VAL	2.2
1	C	1014	GLN	2.1
1	C	939	SER	2.1
1	C	1040	GLN	2.1
1	A	944	PRO	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 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.