



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:32 AM UTC

PDB ID : 1OW7 / pdb_00001ow7
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.; Werner, J.M.; Arold, S.T.
Deposited on : 2003-03-28
Resolution : 2.60 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:

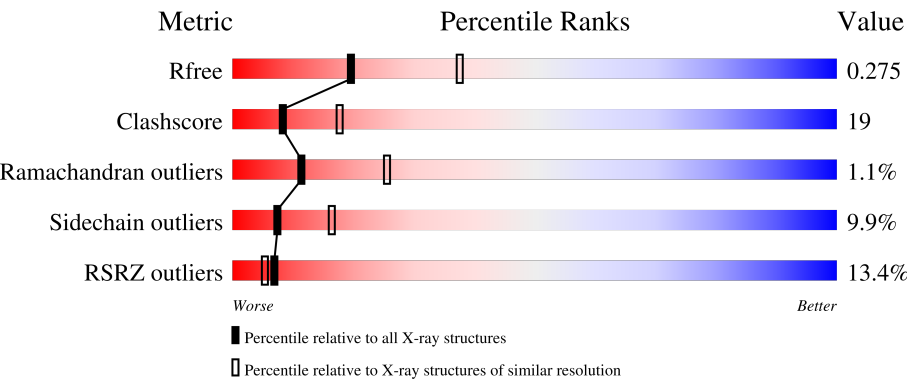
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	5% 51% 26% 6% 17%
1	B	161	4% 58% 25% • 14%
1	C	161	14% 45% 33% 5% • 14%
2	D	13	15% 38% 46% 8% 8%
2	E	13	77% 38% 46% 8% 8%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	F	13	85% 46% 38% 8% 8%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3520 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	E	12	Total	C	N	O	S	0	0	0
			94	56	15	22	1			
2	F	12	Total	C	N	O	S	0	0	0
			94	56	15	22	1			

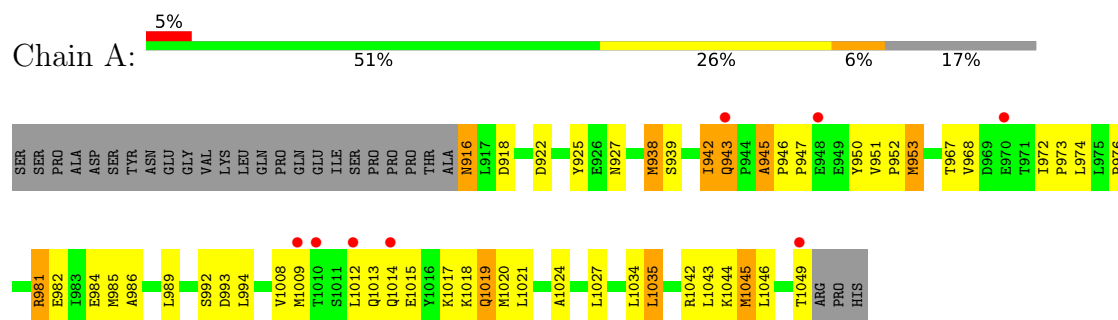
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		
3	B	25	Total	O	0	0
			25	25		
3	C	11	Total	O	0	0
			11	11		
3	D	2	Total	O	0	0
			2	2		
3	F	1	Total	O	0	0
			1	1		

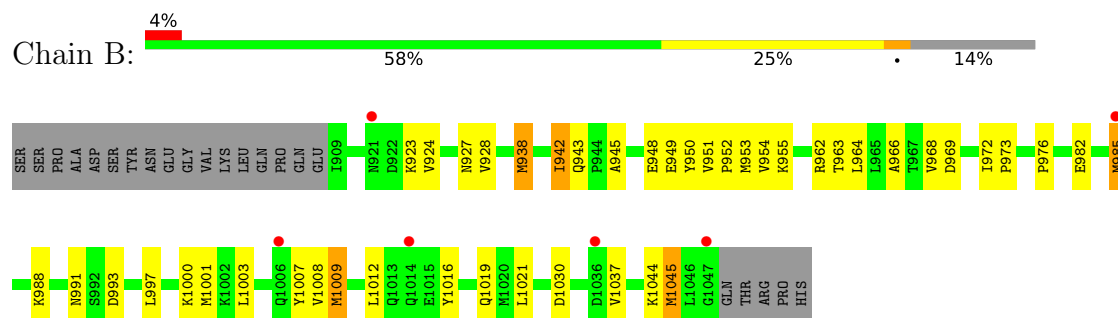
3 Residue-property plots

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.

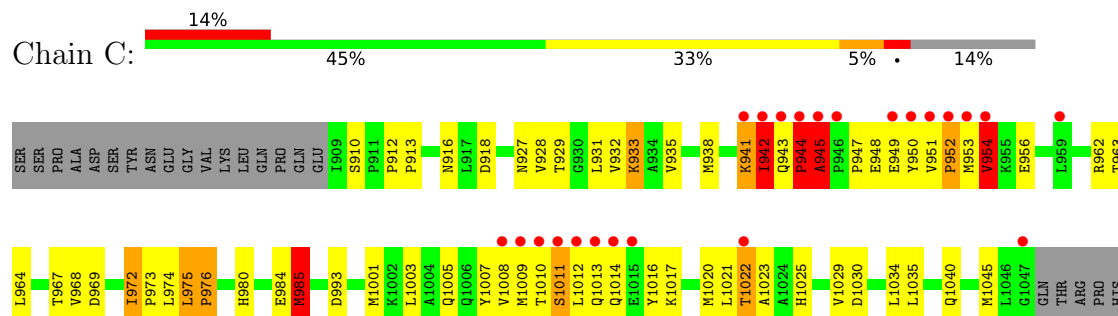
• Molecule 1: Focal adhesion kinase 1



• Molecule 1: Focal adhesion kinase 1

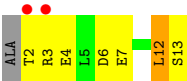


• Molecule 1: Focal adhesion kinase 1

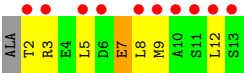
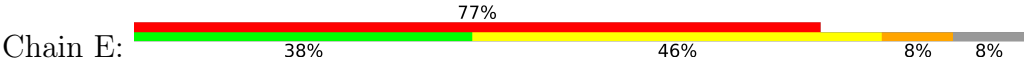


• Molecule 2: Paxillin

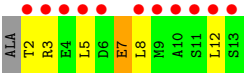
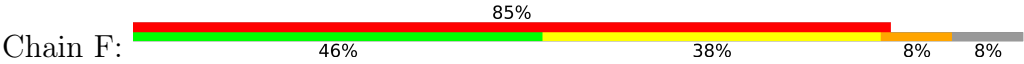




● Molecule 2: Paxillin



● 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	88.99Å 219.66Å 96.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	111.80 – 2.60 109.83 – 2.60	Depositor EDS
% Data completeness (in resolution range)	78.2 (111.80-2.60) 78.1 (109.83-2.60)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.237 , 0.269 0.236 , 0.275	Depositor DCC
R_{free} test set	1189 reflections (5.17%)	wwPDB-VP
Wilson B-factor (Å ²)	63.9	Xtriage
Anisotropy	0.587	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 61.0	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	0.93	EDS
Total number of atoms	3520	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.24% 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	0.72	0/1053	1.11	5/1424 (0.4%)
1	B	0.74	1/1087 (0.1%)	1.13	6/1474 (0.4%)
1	C	0.69	1/1087 (0.1%)	1.17	11/1474 (0.7%)
2	D	0.53	0/93	1.37	2/122 (1.6%)
2	E	0.41	0/93	1.33	1/122 (0.8%)
2	F	0.43	0/93	1.26	1/122 (0.8%)
All	All	0.70	2/3506 (0.1%)	1.15	26/4738 (0.5%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	985	MET	SD-CE	5.18	1.92	1.79
1	B	1009	MET	SD-CE	5.15	1.92	1.79

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	942	ILE	N-CA-C	11.97	121.92	110.42
1	B	942	ILE	N-CA-C	10.79	124.58	111.09
2	F	7	GLU	N-CA-C	-9.51	100.91	111.28
2	E	7	GLU	N-CA-C	-9.20	101.15	111.82
2	D	7	GLU	N-CA-C	-9.16	101.38	111.36
1	A	945	ALA	N-CA-C	8.64	122.85	110.24
1	C	942	ILE	N-CA-C	7.74	125.44	109.34
1	C	954	VAL	N-CA-C	-7.18	104.77	111.45
1	B	1008	VAL	N-CA-C	-7.10	99.44	110.09
1	C	974	LEU	N-CA-C	6.54	119.23	111.71
1	B	924	VAL	N-CA-C	-6.41	104.50	110.53
1	B	928	VAL	N-CA-C	-6.12	104.55	110.42
1	C	972	ILE	CB-CA-C	-6.04	107.95	114.35
1	A	994	LEU	N-CA-C	-5.64	105.21	111.36
1	C	1011	SER	N-CA-C	-5.60	105.95	112.89
1	A	1019	GLN	N-CA-C	-5.47	105.34	112.23

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	928	VAL	N-CA-C	-5.44	105.42	110.53
2	D	3	ARG	N-CA-C	-5.44	105.38	112.23
1	C	945	ALA	CA-C-N	5.34	125.88	120.38
1	C	945	ALA	C-N-CA	5.34	125.88	120.38
1	A	939	SER	N-CA-C	5.21	117.04	111.36
1	C	1040	GLN	N-CA-C	-5.19	105.62	111.28
1	C	945	ALA	N-CA-C	5.14	121.18	109.81
1	C	1020	MET	N-CA-C	-5.13	105.60	111.14
1	B	1007	TYR	N-CA-C	5.07	118.30	112.57
1	B	945	ALA	N-CA-C	5.03	117.02	110.08

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	1042	0	1093	45	0
1	B	1073	0	1127	39	0
1	C	1073	0	1127	47	0
2	D	94	0	92	6	0
2	E	94	0	92	13	0
2	F	94	0	92	6	0
3	A	11	0	0	0	0
3	B	25	0	0	0	0
3	C	11	0	0	0	0
3	D	2	0	0	1	0
3	F	1	0	0	0	0
All	All	3520	0	3623	136	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 19.

All (136) 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:918:ASP:HB3	1:B:1045:MET:CE	2.01	0.91
1:B:985:MET:HG3	1:C:1012:LEU:HG	1.53	0.90
1:B:938:MET:HE2	1:B:954:VAL:HG22	1.57	0.86
1:C:985:MET:HA	1:C:985:MET:HE3	1.60	0.83
1:A:1014:GLN:HE21	1:A:1018:LYS:CE	1.95	0.79
1:A:1014:GLN:HE21	1:A:1018:LYS:HE3	1.47	0.79
1:B:955:LYS:HZ3	2:E:12:LEU:HB2	1.47	0.77
1:C:947:PRO:HB3	1:C:1005:GLN:HE21	1.49	0.77
1:B:962:ARG:HD3	2:E:9:MET:HG3	1.66	0.76
1:A:1008:VAL:HG22	1:A:1009:MET:HG2	1.66	0.75
1:B:991:ASN:OD1	2:E:5:LEU:HD13	1.90	0.72
1:C:1007:TYR:O	1:C:1013:GLN:HB2	1.89	0.71
2:D:4:GLU:HG3	3:D:37:HOH:O	1.89	0.70
1:A:938:MET:HE1	1:A:1024:ALA:HB2	1.72	0.70
1:C:962:ARG:HA	2:F:5:LEU:HD21	1.75	0.69
1:C:938:MET:HE1	1:C:954:VAL:CG2	2.22	0.69
1:C:938:MET:HE1	1:C:954:VAL:HG23	1.75	0.68
1:B:1045:MET:HE3	1:B:1045:MET:HA	1.77	0.67
1:C:950:TYR:HE2	1:C:1017:LYS:HE2	1.60	0.67
1:B:955:LYS:NZ	2:E:12:LEU:HB2	2.10	0.66
1:A:976:PRO:HD3	1:A:1045:MET:HG2	1.78	0.66
1:B:1012:LEU:HG	1:C:985:MET:HG3	1.80	0.64
1:C:968:VAL:O	1:C:972:ILE:HG13	1.98	0.63
1:C:950:TYR:HA	1:C:953:MET:HB2	1.81	0.63
1:B:1001:MET:HE1	2:E:12:LEU:CD2	2.29	0.62
1:C:950:TYR:CE2	1:C:1017:LYS:HE2	2.34	0.62
1:B:938:MET:HE1	1:B:997:LEU:HD11	1.79	0.62
1:C:968:VAL:HG21	1:C:1034:LEU:HD21	1.81	0.62
1:C:985:MET:HA	1:C:985:MET:CE	2.28	0.61
1:C:945:ALA:HB1	1:C:949:GLU:HB2	1.84	0.60
1:A:938:MET:HG2	1:A:953:MET:HE3	1.84	0.60
1:C:1025:HIS:O	1:C:1029:VAL:HG23	2.02	0.60
1:C:942:ILE:HG13	1:C:950:TYR:HB3	1.84	0.59
2:E:3:ARG:O	2:E:7:GLU:HG2	2.02	0.59
1:A:986:ALA:HB3	1:A:1034:LEU:HD13	1.84	0.58
1:C:942:ILE:O	1:C:943:GLN:C	2.45	0.58
1:A:1014:GLN:HE21	1:A:1018:LYS:NZ	2.01	0.58
1:C:943:GLN:HB2	1:C:944:PRO:HD3	1.86	0.57
1:A:1014:GLN:NE2	1:A:1018:LYS:NZ	2.52	0.57
1:A:1009:MET:H	1:A:1013:GLN:HE21	1.53	0.56
1:C:950:TYR:HA	1:C:953:MET:HE3	1.86	0.56
1:C:1010:THR:HG22	1:C:1011:SER:H	1.69	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:943:GLN:CB	1:C:944:PRO:HD3	2.35	0.56
1:B:949:GLU:O	1:B:952:PRO:HD2	2.05	0.56
1:A:1045:MET:HA	1:A:1045:MET:HE2	1.86	0.56
1:A:1014:GLN:NE2	1:A:1018:LYS:CE	2.67	0.56
1:C:927:ASN:ND2	1:C:963:THR:HG22	2.21	0.56
1:A:976:PRO:HD3	1:A:1045:MET:CG	2.36	0.55
1:C:941:LYS:HD3	1:C:941:LYS:C	2.31	0.55
1:C:1003:LEU:HD13	1:C:1016:TYR:CE1	2.41	0.55
1:A:918:ASP:HB3	1:B:1045:MET:HE2	1.86	0.54
1:A:972:ILE:HB	1:A:973:PRO:HD3	1.90	0.54
1:A:986:ALA:CB	1:A:1034:LEU:HD13	2.38	0.53
1:B:1000:LYS:HD2	1:B:1019:GLN:HB3	1.91	0.53
2:D:2:THR:HG22	2:D:2:THR:O	2.08	0.53
1:A:989:LEU:O	1:A:992:SER:HB3	2.09	0.53
1:B:951:VAL:HB	1:B:952:PRO:HD3	1.90	0.53
1:A:916:ASN:OD1	1:A:916:ASN:N	2.42	0.53
1:B:1001:MET:HE1	2:E:12:LEU:HD21	1.89	0.52
1:C:1010:THR:HG22	1:C:1011:SER:N	2.25	0.52
1:B:955:LYS:NZ	2:E:9:MET:O	2.37	0.52
1:B:964:LEU:O	1:B:968:VAL:HG23	2.09	0.52
1:A:1046:LEU:CD1	1:B:1045:MET:HG2	2.40	0.52
1:A:981:ARG:O	1:A:985:MET:HG2	2.10	0.51
1:A:918:ASP:HB3	1:B:1045:MET:HE1	1.87	0.51
1:A:993:ASP:HB2	1:A:1027:LEU:HD13	1.93	0.51
1:C:964:LEU:O	1:C:968:VAL:HG23	2.11	0.50
1:C:976:PRO:HD3	1:C:1045:MET:HB2	1.93	0.50
1:C:927:ASN:ND2	1:C:967:THR:OG1	2.45	0.50
1:B:927:ASN:ND2	1:B:963:THR:HG22	2.27	0.49
1:C:1001:MET:HE1	2:F:12:LEU:HD23	1.93	0.49
1:B:985:MET:CE	1:B:988:LYS:HD2	2.43	0.49
1:C:976:PRO:O	1:C:980:HIS:CD2	2.66	0.49
2:F:3:ARG:O	2:F:7:GLU:HG2	2.12	0.49
1:B:972:ILE:HB	1:B:973:PRO:HD3	1.95	0.49
1:C:972:ILE:N	1:C:973:PRO:CD	2.75	0.49
2:F:5:LEU:O	2:F:8:LEU:HB2	2.13	0.49
1:A:951:VAL:N	1:A:952:PRO:CD	2.75	0.48
2:E:5:LEU:O	2:E:8:LEU:HB2	2.12	0.48
1:B:938:MET:CE	1:B:954:VAL:HG22	2.36	0.48
1:B:1003:LEU:HD13	1:B:1016:TYR:CE1	2.48	0.48
2:F:2:THR:O	2:F:2:THR:HG22	2.14	0.48
1:A:945:ALA:O	1:A:1017:LYS:NZ	2.47	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:947:PRO:HA	1:A:950:TYR:CE2	2.49	0.47
1:B:985:MET:HE3	1:B:988:LYS:HD2	1.96	0.47
1:C:942:ILE:HG23	1:C:943:GLN:N	2.28	0.47
1:C:1022:THR:HG22	1:C:1023:ALA:N	2.29	0.47
1:C:918:ASP:OD1	1:C:918:ASP:C	2.58	0.47
1:A:1035:LEU:HD22	1:A:1035:LEU:O	2.15	0.46
1:A:927:ASN:ND2	1:A:967:THR:OG1	2.49	0.46
1:C:912:PRO:O	1:C:913:PRO:C	2.56	0.46
1:C:976:PRO:CD	1:C:1045:MET:HB2	2.46	0.46
1:B:966:ALA:O	1:B:969:ASP:HB2	2.16	0.46
2:D:2:THR:O	2:D:6:ASP:OD2	2.34	0.45
2:E:2:THR:HG22	2:E:2:THR:O	2.15	0.45
1:A:942:ILE:HG21	1:A:1021:LEU:HD21	1.98	0.45
1:A:942:ILE:HG12	1:A:1017:LYS:HD3	1.98	0.45
1:A:1046:LEU:HD13	1:B:1045:MET:HG2	1.96	0.45
1:C:948:GLU:C	1:C:949:GLU:HG3	2.42	0.45
1:C:972:ILE:HA	1:C:975:LEU:HD12	1.98	0.45
1:B:948:GLU:H	1:B:948:GLU:CD	2.24	0.45
1:A:951:VAL:HB	1:A:952:PRO:HD3	1.98	0.45
1:A:945:ALA:HA	1:A:946:PRO:HD3	1.70	0.44
1:A:1008:VAL:O	1:A:1009:MET:HB2	2.17	0.44
2:F:5:LEU:HA	2:F:8:LEU:HD12	1.99	0.44
1:A:1042:ARG:O	1:A:1046:LEU:HG	2.17	0.44
1:A:1044:LYS:HB3	1:A:1044:LYS:HE2	1.71	0.44
1:C:929:THR:O	1:C:933:LYS:HB2	2.18	0.44
1:B:962:ARG:HG3	2:E:5:LEU:HD21	2.01	0.43
1:A:942:ILE:HG23	1:A:943:GLN:N	2.32	0.43
1:A:976:PRO:HG3	1:A:1045:MET:HE3	2.00	0.43
1:C:993:ASP:OD2	1:C:1030:ASP:OD2	2.35	0.43
2:D:2:THR:HG22	2:D:6:ASP:OD2	2.17	0.43
1:A:1045:MET:HA	1:A:1045:MET:CE	2.48	0.43
1:B:993:ASP:OD2	1:B:1030:ASP:OD2	2.37	0.43
2:D:12:LEU:O	2:D:13:SER:HB2	2.19	0.42
1:A:968:VAL:O	1:A:972:ILE:HG13	2.19	0.42
1:B:1000:LYS:HD2	1:B:1019:GLN:OE1	2.19	0.42
1:C:969:ASP:HA	1:C:972:ILE:HD12	2.00	0.42
1:C:951:VAL:N	1:C:952:PRO:CD	2.82	0.42
2:D:2:THR:O	2:D:2:THR:CG2	2.66	0.42
1:B:976:PRO:HD3	1:B:1045:MET:HB2	2.01	0.42
1:A:1015:GLU:HG3	1:A:1019:GLN:HE21	1.85	0.42
1:B:950:TYR:HA	1:B:953:MET:HE3	2.01	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:951:VAL:HB	1:C:952:PRO:HD3	2.01	0.41
1:A:1043:LEU:HD23	1:A:1043:LEU:HA	1.87	0.41
1:B:942:ILE:HG21	1:B:1021:LEU:HD21	2.02	0.41
1:C:972:ILE:O	1:C:973:PRO:C	2.63	0.41
1:A:974:LEU:HD13	1:B:973:PRO:HB2	2.01	0.41
1:B:1001:MET:HE1	2:E:12:LEU:HD23	2.01	0.41
1:C:938:MET:HE1	1:C:954:VAL:HG22	2.02	0.41
1:A:1008:VAL:O	1:A:1009:MET:CB	2.69	0.41
1:B:951:VAL:N	1:B:952:PRO:CD	2.84	0.41
1:A:922:ASP:HB3	1:A:925:TYR:HB3	2.04	0.40
1:C:931:LEU:O	1:C:932:VAL:C	2.64	0.40
1:B:962:ARG:HG3	2:E:5:LEU:CD2	2.51	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	132/161 (82%)	129 (98%)	3 (2%)	0	100	100
1	B	137/161 (85%)	134 (98%)	3 (2%)	0	100	100
1	C	137/161 (85%)	119 (87%)	13 (10%)	5 (4%)	2	4
2	D	10/13 (77%)	10 (100%)	0	0	100	100
2	E	10/13 (77%)	10 (100%)	0	0	100	100
2	F	10/13 (77%)	10 (100%)	0	0	100	100
All	All	436/522 (84%)	412 (94%)	19 (4%)	5 (1%)	11	25

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	942	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	944	PRO
1	C	916	ASN
1	C	952	PRO
1	C	945	ALA

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%)	105 (90%)	12 (10%)	7	15
1	B	121/141 (86%)	112 (93%)	9 (7%)	13	29
1	C	121/141 (86%)	104 (86%)	17 (14%)	3	6
2	D	11/11 (100%)	10 (91%)	1 (9%)	9	19
2	E	11/11 (100%)	11 (100%)	0	100	100
2	F	11/11 (100%)	11 (100%)	0	100	100
All	All	392/456 (86%)	353 (90%)	39 (10%)	7	16

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	916	ASN
1	A	938	MET
1	A	943	GLN
1	A	953	MET
1	A	981	ARG
1	A	982	GLU
1	A	984	GLU
1	A	1012	LEU
1	A	1020	MET
1	A	1035	LEU
1	A	1045	MET
1	A	1049	THR
1	B	923	LYS
1	B	938	MET

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	943	GLN
1	B	982	GLU
1	B	985	MET
1	B	1009	MET
1	B	1037	VAL
1	B	1044	LYS
1	B	1045	MET
1	C	910	SER
1	C	933	LYS
1	C	935	VAL
1	C	941	LYS
1	C	944	PRO
1	C	954	VAL
1	C	956	GLU
1	C	975	LEU
1	C	976	PRO
1	C	984	GLU
1	C	985	MET
1	C	1008	VAL
1	C	1009	MET
1	C	1014	GLN
1	C	1021	LEU
1	C	1022	THR
1	C	1035	LEU
2	D	12	LEU

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	916	ASN
1	A	927	ASN
1	A	1006	GLN
1	A	1013	GLN
1	A	1014	GLN
1	B	921	ASN
1	C	916	ASN
1	C	927	ASN
1	C	1005	GLN
1	C	1025	HIS
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%)	0.32	8 (5%) 27 22	46, 71, 91, 103	0
1	B	139/161 (86%)	0.12	6 (4%) 40 34	44, 63, 82, 93	0
1	C	139/161 (86%)	0.94	23 (16%) 4 3	49, 81, 127, 136	0
2	D	12/13 (92%)	1.38	2 (16%) 4 3	91, 107, 122, 129	0
2	E	12/13 (92%)	2.93	10 (83%) 0 0	167, 172, 176, 176	0
2	F	12/13 (92%)	2.86	11 (91%) 0 0	163, 169, 176, 178	0
All	All	448/522 (85%)	0.62	60 (13%) 7 5	44, 73, 163, 178	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	951	VAL	5.8
2	F	12	LEU	5.2
2	E	9	MET	5.0
1	C	952	PRO	4.2
1	C	945	ALA	4.1
2	E	8	LEU	3.9
2	E	10	ALA	3.9
1	A	943	GLN	3.8
1	C	1047	GLY	3.8
1	C	1009	MET	3.6
2	E	11	SER	3.6
2	F	2	THR	3.5
1	C	1011	SER	3.4
2	F	6	ASP	3.2
2	E	12	LEU	3.2
1	A	948	GLU	3.2
1	C	1022	THR	3.1
1	A	1009	MET	3.1
1	C	946	PRO	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	F	8	LEU	2.9
2	F	11	SER	2.9
1	B	1014	GLN	2.9
1	C	953	MET	2.9
2	F	10	ALA	2.9
2	E	2	THR	2.9
1	C	944	PRO	2.8
1	A	1014	GLN	2.8
2	F	5	LEU	2.8
2	F	13	SER	2.7
1	A	1010	THR	2.7
2	F	9	MET	2.6
2	E	5	LEU	2.6
2	F	4	GLU	2.6
1	C	954	VAL	2.6
2	E	6	ASP	2.6
1	C	941	LYS	2.6
2	D	2	THR	2.6
1	B	1047	GLY	2.6
1	C	950	TYR	2.6
1	C	949	GLU	2.6
1	C	942	ILE	2.5
1	C	1012	LEU	2.5
1	B	1036	ASP	2.4
1	B	1006	GLN	2.4
1	C	1008	VAL	2.2
1	A	1049	THR	2.2
1	C	943	GLN	2.2
1	C	959	LEU	2.2
2	F	3	ARG	2.1
1	B	921	ASN	2.1
1	A	1012	LEU	2.1
1	B	985	MET	2.1
1	A	970	GLU	2.1
2	D	3	ARG	2.1
2	E	3	ARG	2.1
1	C	1013	GLN	2.1
1	C	1015	GLU	2.0
2	E	13	SER	2.0
1	C	1014	GLN	2.0
1	C	1010	THR	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.