



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:31 PM UTC

PDB ID : 1OW8 / pdb_00001ow8
Title : Paxillin LD2 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 : 2.85 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

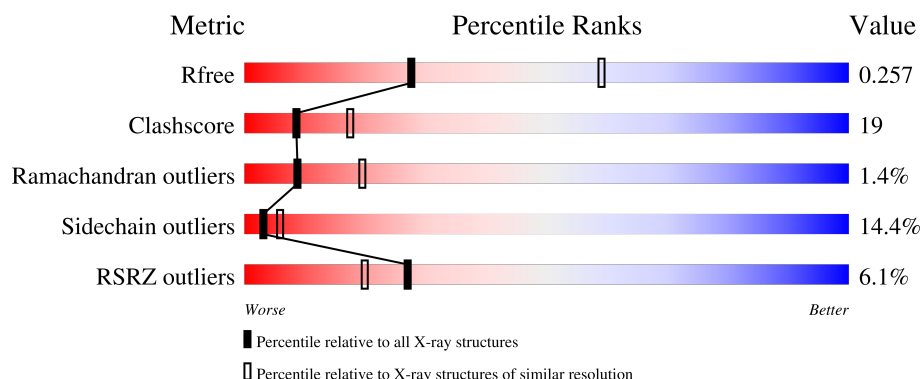
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	15% 48% 29% 6% 17%
1	B	161	52% 30% 6% 12%
1	C	161	6% 49% 31% 7% 12%
2	D	13	15% 31% 38% 23% 8%
2	F	13	85% 62% 23% 8% 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3472 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	141	Total	C	N	O	S	0	0	0
			1091	690	183	211	7			
1	C	142	Total	C	N	O	S	0	0	0
			1098	695	184	212	7			

- Molecule 2 is a protein called Paxillin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	12	Total	C	N	O	0	0	0
			100	63	16	21			
2	F	12	Total	C	N	O	0	0	0
			100	63	16	21			

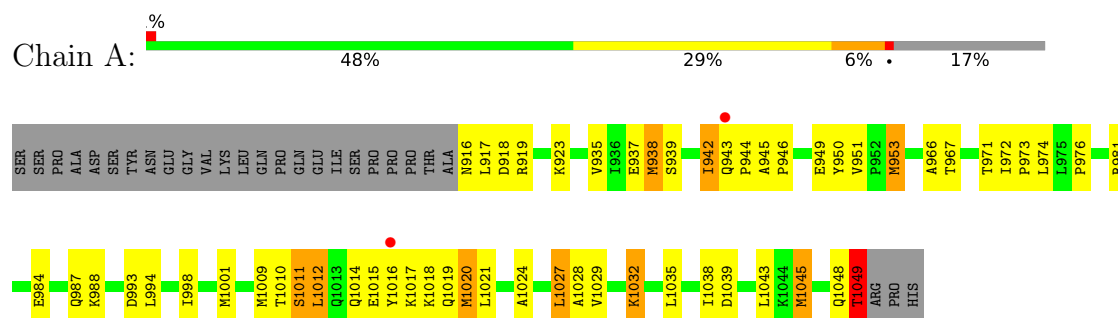
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		
3	B	18	Total	O	0	0
			18	18		
3	C	11	Total	O	0	0
			11	11		
3	D	1	Total	O	0	0
			1	1		

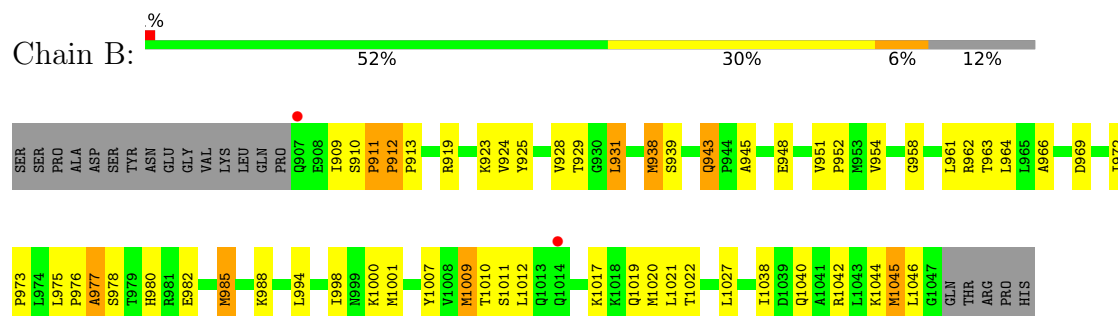
3 Residue-property plots [i](#)

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.

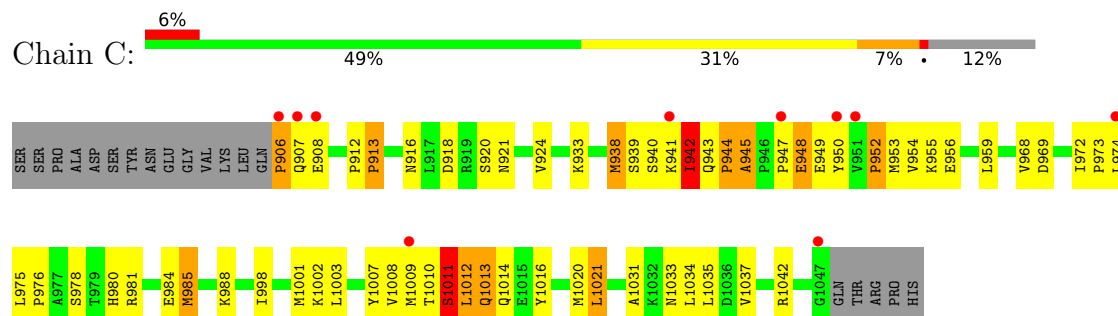
• Molecule 1: Focal adhesion kinase 1



• Molecule 1: Focal adhesion kinase 1



• Molecule 1: Focal adhesion kinase 1

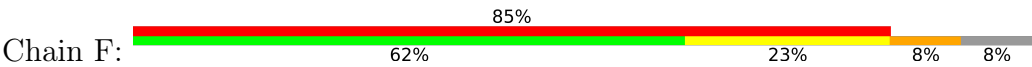


• Molecule 2: Paxillin





● 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	87.98Å 220.78Å 97.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.50 – 2.85 34.50 – 2.85	Depositor EDS
% Data completeness (in resolution range)	99.3 (34.50-2.85) 99.3 (34.50-2.85)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 2.84Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.235 , 0.275 0.224 , 0.257	Depositor DCC
R_{free} test set	1149 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	70.5	Xtriage
Anisotropy	0.445	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3472	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% 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	0.71	1/1053 (0.1%)	1.44	7/1424 (0.5%)
1	B	0.72	1/1105 (0.1%)	1.44	7/1498 (0.5%)
1	C	0.69	0/1113	1.31	14/1509 (0.9%)
2	D	0.85	0/99	1.42	3/131 (2.3%)
2	F	0.72	0/99	1.18	1/131 (0.8%)
All	All	0.71	2/3469 (0.1%)	1.39	32/4693 (0.7%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1009	MET	SD-CE	6.14	1.95	1.79
1	A	1009	MET	SD-CE	5.02	1.92	1.79

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	910	SER	CA-C-N	10.87	127.59	119.66
1	B	910	SER	C-N-CA	10.87	127.59	119.66
1	B	931	LEU	N-CA-C	-9.56	101.17	113.12
1	B	911	PRO	CB-CA-C	-8.82	103.09	111.39
1	C	907	GLN	N-CA-C	8.69	121.52	108.31
1	A	994	LEU	N-CA-C	-8.61	101.83	111.82
1	C	945	ALA	CA-C-N	8.41	129.05	120.38
1	C	945	ALA	C-N-CA	8.41	129.05	120.38
1	A	1049	THR	CA-C-O	8.23	134.80	120.80
1	C	942	ILE	N-CA-C	7.78	125.52	109.34
2	D	3	SER	N-CA-C	-7.66	102.90	111.71
1	C	945	ALA	N-CA-C	7.13	125.56	109.81
1	A	937	GLU	N-CA-C	-6.97	103.68	111.28
1	A	917	LEU	N-CA-C	6.83	119.87	110.24
1	C	907	GLN	CB-CA-C	-6.48	109.09	116.54
1	C	1011	SER	N-CA-C	-6.41	104.39	111.82
1	C	906	PRO	CA-C-N	6.02	133.25	122.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	906	PRO	C-N-CA	6.02	133.25	122.89
1	C	924	VAL	N-CA-C	-5.97	104.91	110.53
1	C	1031	ALA	N-CA-C	-5.96	104.91	111.82
2	D	8	LEU	N-CA-C	-5.87	101.94	111.04
1	A	987	GLN	N-CA-C	-5.67	105.17	111.36
1	B	911	PRO	CA-C-O	5.52	124.88	120.90
1	C	1001	MET	N-CA-C	-5.50	103.14	111.56
1	B	912	PRO	N-CA-C	-5.31	105.37	110.47
1	A	1027	LEU	N-CA-C	-5.30	105.14	111.03
2	D	5	LEU	N-CA-C	-5.27	105.54	111.28
1	A	939	SER	N-CA-C	-5.25	105.63	111.36
1	B	977	ALA	N-CA-C	-5.25	106.13	112.54
1	C	959	LEU	N-CA-C	-5.14	105.13	111.40
1	C	1042	ARG	N-CA-C	-5.11	105.41	111.69
2	F	8	LEU	N-CA-C	-5.04	105.43	111.03

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	41	0
1	B	1091	0	1141	35	0
1	C	1098	0	1149	53	0
2	D	100	0	105	10	0
2	F	100	0	105	7	0
3	A	11	0	0	0	0
3	B	18	0	0	2	0
3	C	11	0	0	1	0
3	D	1	0	0	0	0
All	All	3472	0	3593	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 19.

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:C:985:MET:HA	1:C:985:MET:HE3	1.50	0.91
2:D:5:LEU:HA	2:D:8:LEU:HD12	1.54	0.89
1:B:1045:MET:HE3	1:B:1045:MET:HA	1.61	0.83
1:C:985:MET:HA	1:C:985:MET:CE	2.09	0.82
1:C:955:LYS:HA	2:F:12:LEU:HD13	1.62	0.81
2:F:5:LEU:HA	2:F:8:LEU:HD12	1.62	0.80
1:A:1014:GLN:HE21	1:A:1018:LYS:NZ	1.82	0.78
1:A:918:ASP:HB3	1:B:1045:MET:CE	2.16	0.76
1:C:942:ILE:O	1:C:943:GLN:C	2.29	0.75
1:B:978:SER:HB3	3:B:35:HOH:O	1.90	0.71
1:B:985:MET:HG3	1:C:1012:LEU:HG	1.75	0.68
1:C:985:MET:HE3	1:C:985:MET:CA	2.23	0.66
1:C:969:ASP:HA	1:C:972:ILE:HD12	1.78	0.66
1:C:945:ALA:HB1	1:C:949:GLU:HB2	1.77	0.66
2:F:2:LEU:HD11	2:F:5:LEU:HD23	1.78	0.66
1:A:1011:SER:C	1:A:1012:LEU:HD23	2.23	0.64
1:C:943:GLN:HB2	1:C:944:PRO:CD	2.27	0.64
2:D:2:LEU:HD11	2:D:5:LEU:HD23	1.81	0.63
1:C:955:LYS:CA	2:F:12:LEU:HD13	2.28	0.63
1:C:943:GLN:CB	1:C:944:PRO:CD	2.77	0.63
1:C:976:PRO:O	1:C:980:HIS:CD2	2.54	0.61
1:A:945:ALA:O	1:A:1017:LYS:NZ	2.33	0.61
1:B:966:ALA:O	1:B:969:ASP:HB2	1.99	0.61
1:A:974:LEU:HD13	1:B:973:PRO:HB2	1.81	0.61
1:C:943:GLN:HB2	1:C:944:PRO:HD2	1.83	0.61
1:C:947:PRO:HG3	1:C:950:TYR:OH	2.01	0.60
1:B:1011:SER:OG	1:C:985:MET:HG2	2.02	0.59
1:C:972:ILE:N	1:C:973:PRO:CD	2.65	0.59
1:A:1032:LYS:HG3	2:D:9:LEU:HD11	1.83	0.59
1:B:958:GLY:C	1:B:962:ARG:NH1	2.62	0.58
1:B:1012:LEU:HG	1:C:985:MET:HG3	1.85	0.58
1:A:918:ASP:HB3	1:B:1045:MET:HE2	1.85	0.57
1:A:943:GLN:HB3	1:A:944:PRO:HD3	1.86	0.57
1:C:972:ILE:HA	1:C:975:LEU:HD12	1.86	0.56
1:A:1012:LEU:HD23	1:A:1012:LEU:N	2.21	0.56
1:A:1048:GLN:HG2	1:A:1049:THR:H	1.68	0.56
1:A:972:ILE:HB	1:A:973:PRO:HD3	1.88	0.56
1:C:921:ASN:HB2	3:C:53:HOH:O	2.06	0.56
2:D:2:LEU:HD12	2:D:5:LEU:HB3	1.88	0.55
1:B:951:VAL:HB	1:B:952:PRO:HD3	1.89	0.55
2:F:2:LEU:HD12	2:F:5:LEU:HB3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:938:MET:HG2	1:A:953:MET:HE3	1.89	0.54
1:A:1014:GLN:HE21	1:A:1018:LYS:CE	2.21	0.54
1:A:935:VAL:HG13	1:A:1024:ALA:HB1	1.90	0.53
1:A:1032:LYS:O	1:A:1032:LYS:HD3	2.08	0.53
2:F:2:LEU:CD1	2:F:5:LEU:HD23	2.38	0.53
1:A:946:PRO:HG2	1:A:949:GLU:HG3	1.90	0.53
1:B:1000:LYS:HD2	1:B:1019:GLN:CB	2.39	0.53
1:B:1000:LYS:HD2	1:B:1019:GLN:HB3	1.90	0.53
1:A:938:MET:HG3	1:A:953:MET:HB3	1.91	0.53
1:A:1014:GLN:HE21	1:A:1018:LYS:HZ2	1.54	0.53
2:D:2:LEU:CD1	2:D:5:LEU:HD23	2.40	0.52
1:A:1014:GLN:NE2	1:A:1018:LYS:NZ	2.55	0.52
1:C:952:PRO:C	1:C:954:VAL:N	2.66	0.52
1:A:919:ARG:NH2	1:A:1039:ASP:OD2	2.37	0.52
1:C:942:ILE:HG23	1:C:943:GLN:N	2.24	0.51
1:C:968:VAL:HG21	1:C:1034:LEU:HD21	1.91	0.51
1:B:911:PRO:O	1:B:912:PRO:C	2.52	0.51
1:A:951:VAL:HG22	1:A:1001:MET:SD	2.51	0.51
1:A:1032:LYS:HG3	2:D:9:LEU:CD1	2.41	0.51
1:C:950:TYR:HA	1:C:953:MET:HB2	1.93	0.51
1:A:1032:LYS:CG	2:D:9:LEU:HD11	2.41	0.50
1:C:1010:THR:HG22	1:C:1011:SER:H	1.75	0.50
1:A:942:ILE:HG21	1:A:1021:LEU:HD21	1.94	0.50
1:C:938:MET:HG2	1:C:939:SER:N	2.25	0.50
1:C:968:VAL:HG12	1:C:972:ILE:HD11	1.93	0.49
1:A:993:ASP:HB2	1:A:1027:LEU:HD13	1.95	0.49
1:A:1020:MET:O	1:A:1021:LEU:C	2.56	0.49
1:C:972:ILE:N	1:C:973:PRO:HD3	2.28	0.48
1:B:938:MET:HE2	1:B:954:VAL:HG22	1.95	0.48
1:B:963:THR:O	1:B:964:LEU:C	2.55	0.48
1:C:938:MET:HE1	1:C:954:VAL:HG22	1.93	0.48
1:A:1015:GLU:HG3	1:A:1019:GLN:HE21	1.79	0.48
1:B:985:MET:HE3	1:B:988:LYS:HD2	1.95	0.48
1:C:943:GLN:CB	1:C:944:PRO:HD3	2.43	0.48
1:A:976:PRO:HD3	1:A:1045:MET:HG2	1.96	0.48
1:A:1014:GLN:C	1:A:1016:TYR:N	2.70	0.47
1:B:919:ARG:HD2	1:B:925:TYR:CD2	2.49	0.47
2:D:8:LEU:O	2:D:9:LEU:C	2.57	0.47
1:A:1048:GLN:HG2	1:A:1049:THR:N	2.29	0.47
1:C:972:ILE:HG22	1:C:980:HIS:CE1	2.49	0.47
2:D:6:ASP:HA	2:D:9:LEU:HD12	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:942:ILE:HG21	1:C:1021:LEU:HD21	1.97	0.47
1:B:943:GLN:O	1:B:1017:LYS:NZ	2.46	0.47
1:A:976:PRO:HD3	1:A:1045:MET:CG	2.44	0.47
1:B:939:SER:CB	1:B:1021:LEU:HD22	2.44	0.47
1:C:955:LYS:HA	2:F:12:LEU:CD1	2.38	0.47
1:B:1038:ILE:O	1:B:1038:ILE:HG22	2.13	0.47
1:C:939:SER:HA	1:C:1021:LEU:HD22	1.97	0.47
1:C:948:GLU:C	1:C:949:GLU:HG3	2.40	0.46
1:B:924:VAL:O	1:B:928:VAL:HG23	2.16	0.46
1:B:945:ALA:O	1:B:1017:LYS:NZ	2.37	0.46
1:C:1003:LEU:HD13	1:C:1016:TYR:CE1	2.51	0.46
1:B:1042:ARG:O	1:B:1046:LEU:HG	2.16	0.46
2:D:8:LEU:O	2:D:11:GLU:N	2.49	0.46
1:A:950:TYR:HA	1:A:953:MET:HE2	1.98	0.45
1:B:912:PRO:O	1:B:913:PRO:C	2.60	0.45
1:C:968:VAL:O	1:C:972:ILE:CD1	2.65	0.45
1:C:998:ILE:CG2	1:C:1002:LYS:HE2	2.46	0.45
1:C:1010:THR:HG22	1:C:1011:SER:N	2.31	0.45
1:C:952:PRO:O	1:C:953:MET:C	2.61	0.45
1:B:985:MET:CE	1:B:988:LYS:HD2	2.48	0.44
1:C:1033:ASN:O	1:C:1037:VAL:HG23	2.18	0.44
1:B:972:ILE:HB	1:B:973:PRO:HD3	1.99	0.44
1:C:942:ILE:CG2	1:C:943:GLN:N	2.80	0.44
1:C:1007:TYR:O	1:C:1013:GLN:HB2	2.18	0.44
1:C:918:ASP:CG	1:C:920:SER:HG	2.25	0.44
1:A:1027:LEU:O	1:A:1028:ALA:C	2.60	0.43
1:A:1045:MET:HA	1:A:1045:MET:CE	2.48	0.43
1:C:952:PRO:O	1:C:954:VAL:N	2.51	0.43
1:A:971:THR:HG22	1:A:971:THR:O	2.17	0.43
1:C:976:PRO:O	1:C:980:HIS:HD2	1.98	0.43
1:B:975:LEU:O	1:B:976:PRO:C	2.61	0.43
1:C:968:VAL:O	1:C:972:ILE:HD12	2.18	0.43
1:B:994:LEU:O	1:B:998:ILE:HG13	2.17	0.43
1:B:1045:MET:HA	1:B:1045:MET:CE	2.41	0.43
1:C:1020:MET:O	1:C:1021:LEU:C	2.61	0.42
1:B:977:ALA:HA	1:B:980:HIS:CD2	2.55	0.42
1:B:1001:MET:HA	1:B:1020:MET:SD	2.59	0.42
1:A:972:ILE:N	1:A:973:PRO:CD	2.82	0.42
1:C:998:ILE:HG22	1:C:1002:LYS:HE2	2.02	0.42
1:C:969:ASP:O	1:C:973:PRO:HD3	2.20	0.42
1:A:1043:LEU:HD23	1:A:1043:LEU:HA	1.80	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:998:ILE:HG22	1:C:1002:LYS:CE	2.50	0.42
1:C:952:PRO:O	1:C:955:LYS:N	2.53	0.41
1:A:1014:GLN:C	1:A:1016:TYR:H	2.29	0.41
1:B:909:ILE:HD12	3:B:19:HOH:O	2.21	0.41
1:B:1007:TYR:OH	1:C:988:LYS:HG2	2.21	0.41
1:C:912:PRO:O	1:C:913:PRO:C	2.64	0.41
1:A:1014:GLN:HE21	1:A:1018:LYS:HE3	1.85	0.41
1:A:1014:GLN:O	1:A:1015:GLU:C	2.60	0.41
1:B:961:LEU:O	1:B:961:LEU:HD12	2.21	0.41
1:A:966:ALA:O	1:A:967:THR:C	2.63	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	31
1	B	139/161 (86%)	133 (96%)	6 (4%)	0	100	100
1	C	140/161 (87%)	120 (86%)	15 (11%)	5 (4%)	2	5
2	D	10/13 (77%)	8 (80%)	2 (20%)	0	100	100
2	F	10/13 (77%)	10 (100%)	0	0	100	100
All	All	431/509 (85%)	398 (92%)	27 (6%)	6 (1%)	9	19

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	916	ASN
1	C	942	ILE
1	C	908	GLU
1	A	1011	SER

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Mol	Chain	Res	Type
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%)	99 (85%)	18 (15%)	2	5
1	B	123/141 (87%)	108 (88%)	15 (12%)	5	9
1	C	124/141 (88%)	103 (83%)	21 (17%)	2	3
2	D	12/13 (92%)	10 (83%)	2 (17%)	2	3
2	F	12/13 (92%)	12 (100%)	0	100	100
All	All	388/449 (86%)	332 (86%)	56 (14%)	3	6

All (56) 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	942	ILE
1	A	953	MET
1	A	981	ARG
1	A	984	GLU
1	A	988	LYS
1	A	998	ILE
1	A	1010	THR
1	A	1012	LEU
1	A	1020	MET
1	A	1029	VAL
1	A	1032	LYS
1	A	1035	LEU
1	A	1038	ILE
1	A	1045	MET

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Mol	Chain	Res	Type
1	A	1049	THR
1	B	923	LYS
1	B	929	THR
1	B	931	LEU
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	1010	THR
1	B	1022	THR
1	B	1027	LEU
1	B	1040	GLN
1	B	1044	LYS
1	B	1045	MET
1	C	906	PRO
1	C	913	PRO
1	C	933	LYS
1	C	938	MET
1	C	940	SER
1	C	941	LYS
1	C	948	GLU
1	C	956	GLU
1	C	974	LEU
1	C	978	SER
1	C	981	ARG
1	C	984	GLU
1	C	985	MET
1	C	1008	VAL
1	C	1009	MET
1	C	1011	SER
1	C	1012	LEU
1	C	1013	GLN
1	C	1014	GLN
1	C	1021	LEU
1	C	1035	LEU
2	D	3	SER
2	D	4	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	916	ASN
1	A	987	GLN
1	A	1005	GLN
1	A	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%)	-0.07	2 (1%) 72 65	41, 64, 90, 104	0
1	B	141/161 (87%)	-0.31	2 (1%) 73 67	40, 59, 76, 123	0
1	C	142/161 (88%)	0.37	10 (7%) 22 16	51, 76, 120, 132	0
2	D	12/13 (92%)	1.31	2 (16%) 4 3	83, 90, 112, 118	0
2	F	12/13 (92%)	4.43	11 (91%) 0 0	118, 134, 138, 147	0
All	All	441/509 (86%)	0.15	27 (6%) 27 20	40, 68, 119, 147	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	12	LEU	7.3
2	F	9	LEU	6.8
2	F	2	LEU	6.3
2	F	8	LEU	6.3
2	F	10	LEU	5.9
1	C	906	PRO	5.2
2	F	5	LEU	4.9
2	D	2	LEU	4.2
1	C	1009	MET	3.6
1	B	907	GLN	3.6
2	F	3	SER	3.4
2	F	13	ASN	3.0
1	C	941	LYS	3.0
1	C	907	GLN	2.8
2	F	6	ASP	2.8
1	A	1016	TYR	2.7
1	C	1047	GLY	2.6
2	F	4	GLU	2.6
2	F	7	ARG	2.4
1	C	950	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
2	D	3	SER	2.4
1	C	947	PRO	2.3
1	C	951	VAL	2.3
1	C	974	LEU	2.3
1	C	908	GLU	2.2
1	A	943	GLN	2.2
1	B	1014	GLN	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.