



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:32 AM GMT

PDB ID : 3B71
Title : CD4 endocytosis motif bound to the Focal Adhesion Targeting (FAT) domain
of the Focal Adhesion Kinase
Authors : Garron, M.-L.; Arold, S.T.
Deposited on : 2007-10-30
Resolution : 2.82 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

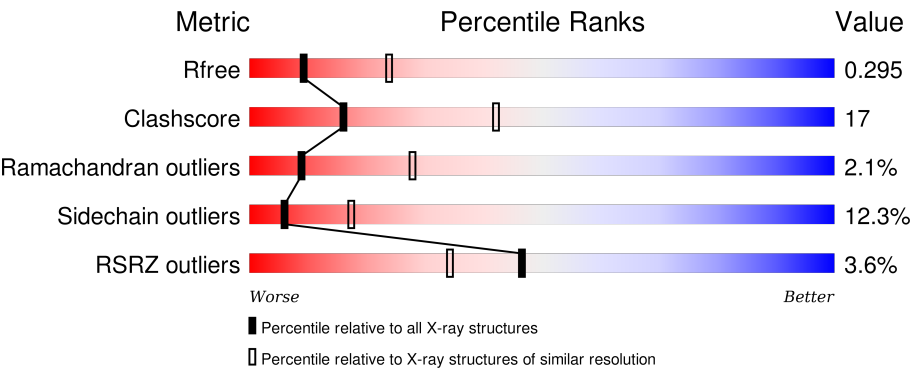
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.82 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}	91344	2676 (2.84-2.80)
Clashscore	102246	3124 (2.84-2.80)
Ramachandran outliers	100387	3072 (2.84-2.80)
Sidechain outliers	100360	3074 (2.84-2.80)
RSRZ outliers	91569	2690 (2.84-2.80)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	162	
1	B	162	
1	C	162	
2	D	23	
2	E	23	

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Mol	Chain	Length	Quality of chain
2	F	23	<div><div><div></div><div></div><div></div><div></div></div><div>4%13%17%9%61%</div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3454 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.

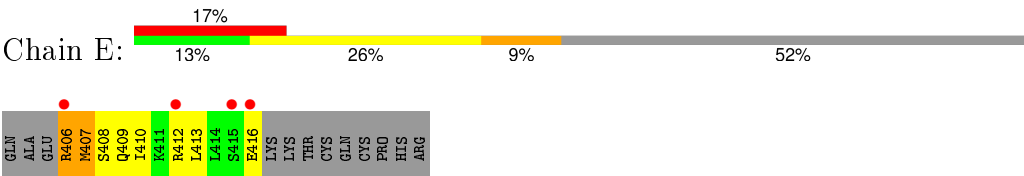
- Molecule 1 is a protein called Focal adhesion kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	134	Total	C	N	O	S	0	1	0
			1048	662	178	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			

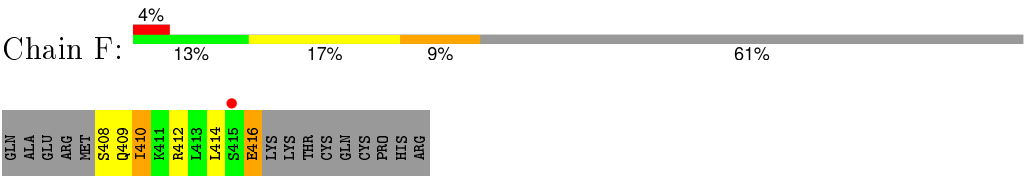
- Molecule 2 is a protein called T-cell surface glycoprotein CD4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	11	Total	C	N	O	S	0	0	0
			93	57	19	16	1			
2	E	11	Total	C	N	O	S	0	0	0
			93	57	19	16	1			
2	F	9	Total	C	N	O		0	0	0
			74	46	14	14				

● Molecule 2: T-cell surface glycoprotein CD4



● Molecule 2: T-cell surface glycoprotein CD4



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	88.07Å 220.47Å 97.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.00 – 2.82 29.35 – 2.82	Depositor EDS
% Data completeness (in resolution range)	93.8 (29.00-2.82) 93.8 (29.35-2.82)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.33 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.241 , 0.307 0.233 , 0.295	Depositor DCC
R_{free} test set	459 reflections (2.14%)	DCC
Wilson B-factor (Å ²)	69.3	Xtriage
Anisotropy	0.214	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 39.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 21922 reflections (0.005%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3454	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.66% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.86	0/1064	0.90	1/1439 (0.1%)
1	B	0.87	0/1087	0.90	0/1474
1	C	0.73	0/1087	0.89	1/1474 (0.1%)
2	D	0.54	0/92	0.92	0/119
2	E	0.78	0/92	0.74	0/119
2	F	1.05	0/73	0.74	0/95
All	All	0.82	0/3495	0.89	2/4720 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	917	LEU	CB-CG-CD1	-5.76	101.20	111.00
1	C	962	ARG	NE-CZ-NH1	5.28	122.94	120.30

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	1048	0	1096	23	0
1	B	1073	0	1127	41	0
1	C	1073	0	1127	38	0
2	D	93	0	104	10	0
2	E	93	0	104	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	74	0	82	4	0
All	All	3454	0	3640	120	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 17.

All (120) 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:942:ILE:HG23	1:C:943:GLN:N	1.82	0.94
1:B:950:TYR:HA	1:B:953:MET:HE2	1.51	0.92
1:C:977:ALA:HA	1:C:980:HIS:NE2	1.84	0.91
1:C:1003:LEU:O	1:C:1006:GLN:HB2	1.71	0.91
1:B:950:TYR:HA	1:B:953:MET:CE	2.00	0.90
1:B:1008:VAL:HA	1:B:1013:GLN:OE1	1.74	0.85
1:C:942:ILE:CG2	1:C:943:GLN:H	1.90	0.83
1:C:942:ILE:HG23	1:C:943:GLN:H	1.37	0.83
2:F:412:ARG:O	2:F:416:GLU:HB3	1.82	0.79
1:B:956:GLU:OE2	1:B:956:GLU:HA	1.82	0.79
1:C:942:ILE:CG2	1:C:943:GLN:N	2.47	0.78
1:B:943:GLN:HB3	1:B:944:PRO:HD3	1.66	0.78
1:C:977:ALA:HA	1:C:980:HIS:CD2	2.19	0.77
1:C:968:VAL:O	1:C:972:ILE:HG13	1.86	0.76
2:E:406:ARG:O	2:E:410:ILE:HG22	1.86	0.76
2:D:406:ARG:HB3	2:D:406:ARG:HH11	1.49	0.76
2:D:413:LEU:HA	2:D:416:GLU:CD	2.06	0.75
1:C:980:HIS:O	1:C:984:GLU:HG2	1.92	0.70
2:E:408:SER:OG	2:E:412:ARG:NH1	2.24	0.69
1:B:995:GLY:HA3	2:E:406:ARG:HH21	1.58	0.69
1:A:1043:LEU:HD13	1:A:1048:GLN:HG2	1.76	0.68
1:A:998:ILE:CG2	1:A:1002:LYS:HE3	2.25	0.67
1:A:998:ILE:HG22	1:A:1002:LYS:HE3	1.78	0.66
1:A:942:ILE:HG21	1:A:1021:LEU:HD11	1.80	0.64
1:A:1025[A]:HIS:CE1	2:D:406:ARG:HH12	2.16	0.64
1:B:950:TYR:HA	1:B:953:MET:HE3	1.79	0.64
1:C:977:ALA:HA	1:C:980:HIS:HE2	1.62	0.63
1:B:979:THR:O	1:B:983:ILE:HD12	1.99	0.63
1:C:1007:TYR:O	1:C:1013:GLN:HB3	1.98	0.63
1:B:937:GLU:HA	1:B:937:GLU:OE1	1.99	0.62
1:C:962:ARG:HG3	1:C:962:ARG:HH11	1.63	0.62
1:B:976:PRO:HB2	1:B:979:THR:HG23	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:938:MET:O	1:C:942:ILE:HB	2.01	0.60
2:D:414:LEU:O	2:D:414:LEU:HD22	2.01	0.60
2:D:412:ARG:NH1	2:D:412:ARG:HB2	2.17	0.60
1:C:938:MET:HE1	1:C:954:VAL:HG22	1.82	0.60
1:C:951:VAL:HB	1:C:952:PRO:HD3	1.84	0.59
1:B:998:ILE:HD13	2:E:409:GLN:HB3	1.84	0.59
1:B:950:TYR:HE2	1:B:1013:GLN:NE2	1.99	0.59
1:B:950:TYR:CA	1:B:953:MET:HE2	2.28	0.59
1:C:969:ASP:O	1:C:973:PRO:HD3	2.04	0.58
2:D:409:GLN:HA	2:D:412:ARG:NH2	2.19	0.58
1:A:1012:LEU:O	1:A:1013:GLN:C	2.42	0.57
1:B:949:GLU:O	1:B:952:PRO:HD2	2.05	0.57
1:C:979:THR:HG23	1:C:979:THR:O	2.05	0.56
1:C:910:SER:HB2	1:C:911:PRO:HD2	1.88	0.55
1:B:985:MET:HE2	1:B:988:LYS:HD3	1.88	0.55
1:B:980:HIS:O	1:B:984:GLU:HG3	2.06	0.55
1:B:949:GLU:C	1:B:952:PRO:HD2	2.27	0.55
2:F:408:SER:N	2:F:410:ILE:HG23	2.22	0.55
1:C:950:TYR:O	1:C:954:VAL:HG23	2.07	0.55
1:C:950:TYR:HA	1:C:953:MET:HB2	1.89	0.54
2:D:412:ARG:HH11	2:D:412:ARG:HB2	1.71	0.54
1:B:931:LEU:O	1:B:931:LEU:HD12	2.07	0.54
1:A:943:GLN:OE1	1:A:1014:GLN:HG3	2.08	0.53
1:B:943:GLN:HB3	1:B:944:PRO:CD	2.37	0.53
1:C:982:GLU:HG2	1:C:1037:VAL:HG11	1.91	0.52
1:A:1033:ASN:O	1:A:1037:VAL:HG23	2.10	0.51
2:E:407:MET:HA	2:E:410:ILE:CG2	2.41	0.51
1:B:1008:VAL:O	1:B:1008:VAL:HG22	2.10	0.51
2:F:409:GLN:HA	2:F:412:ARG:HB2	1.94	0.50
1:B:1018:LYS:HA	1:B:1021:LEU:HD12	1.94	0.49
1:A:972:ILE:HB	1:A:973:PRO:HD3	1.95	0.49
1:B:919:ARG:HD2	1:B:925:TYR:CD2	2.48	0.49
1:C:977:ALA:O	1:C:979:THR:N	2.46	0.48
1:C:977:ALA:C	1:C:979:THR:H	2.16	0.48
1:A:961:LEU:HD23	1:A:994:LEU:HD22	1.95	0.48
1:B:950:TYR:CA	1:B:953:MET:CE	2.85	0.48
2:E:406:ARG:N	2:E:406:ARG:HE	2.12	0.48
2:E:406:ARG:C	2:E:410:ILE:HG22	2.33	0.48
1:A:937:GLU:OE2	1:A:941:LYS:HE3	2.13	0.48
1:A:971:THR:HG22	1:A:975:LEU:HD11	1.96	0.48
2:F:410:ILE:HG13	2:F:414:LEU:HD12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:942:ILE:O	1:B:943:GLN:C	2.52	0.47
1:C:931:LEU:HD23	1:C:1031:ALA:HB2	1.96	0.47
1:B:1003:LEU:O	1:B:1007:TYR:HD1	1.97	0.47
1:A:993:ASP:CB	1:A:1027:LEU:HD13	2.45	0.47
1:A:986:ALA:HB3	1:A:1034:LEU:HD13	1.96	0.47
1:B:909:ILE:HG23	1:B:909:ILE:O	2.15	0.46
1:A:971:THR:O	1:A:972:ILE:C	2.54	0.45
1:B:990:LEU:HD22	1:B:1027:LEU:HD12	1.98	0.45
1:B:950:TYR:CE2	1:B:1013:GLN:NE2	2.81	0.45
1:B:994:LEU:HD23	2:E:410:ILE:HB	1.98	0.45
1:C:990:LEU:HD12	1:C:990:LEU:N	2.32	0.45
1:C:983:ILE:O	1:C:984:GLU:C	2.55	0.45
1:A:993:ASP:HB2	1:A:1027:LEU:CD1	2.47	0.45
1:A:965:LEU:HD22	1:A:987:GLN:HG2	1.99	0.45
1:C:942:ILE:HG21	1:C:1021:LEU:HD21	1.99	0.45
1:A:976:PRO:HD3	1:A:1045:MET:SD	2.57	0.45
1:B:1034:LEU:O	1:B:1035:LEU:C	2.54	0.45
1:C:964:LEU:O	1:C:968:VAL:HG23	2.17	0.45
1:C:972:ILE:HG22	1:C:980:HIS:ND1	2.33	0.44
2:E:407:MET:CA	2:E:410:ILE:HG22	2.47	0.44
1:A:918:ASP:HB3	1:B:1045:MET:SD	2.58	0.44
1:A:1042:ARG:O	1:A:1046:LEU:N	2.46	0.44
1:A:976:PRO:HD3	1:A:1045:MET:HB2	2.00	0.44
1:C:1034:LEU:O	1:C:1037:VAL:HG23	2.18	0.44
1:B:961:LEU:O	1:B:961:LEU:HD12	2.18	0.43
1:C:972:ILE:N	1:C:973:PRO:HD2	2.33	0.42
1:B:1001:MET:O	1:B:1005:GLN:HG3	2.19	0.42
1:B:910:SER:HB2	1:B:911:PRO:HD2	2.00	0.42
1:A:1038:ILE:O	1:A:1041:ALA:HB3	2.19	0.42
1:C:971:THR:C	1:C:973:PRO:HD2	2.40	0.42
1:C:969:ASP:O	1:C:973:PRO:CD	2.67	0.42
1:B:959:LEU:HA	1:B:959:LEU:HD23	1.92	0.42
1:C:1013:GLN:HG3	1:C:1014:GLN:N	2.35	0.42
1:C:1037:VAL:O	1:C:1038:ILE:C	2.57	0.42
1:B:929:THR:O	1:B:933:LYS:HG2	2.19	0.42
1:C:1045:MET:O	1:C:1045:MET:HG2	2.20	0.42
2:D:410:ILE:H	2:D:410:ILE:HG13	1.74	0.41
2:D:412:ARG:HH11	2:D:412:ARG:CB	2.32	0.41
1:C:981:ARG:HB3	1:C:981:ARG:HE	1.61	0.41
1:B:1010:THR:HB	1:C:981:ARG:HH22	1.85	0.41
1:B:945:ALA:HA	1:B:946:PRO:HD3	1.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:409:GLN:O	2:E:413:LEU:N	2.32	0.40
1:A:916:ASN:HB3	1:A:917:LEU:H	1.63	0.40
1:B:1044:LYS:HE2	1:B:1044:LYS:HB3	1.81	0.40
1:B:942:ILE:HG21	1:B:1021:LEU:HD21	2.03	0.40
1:B:948:GLU:HG3	1:B:948:GLU:H	1.41	0.40
2:D:406:ARG:CB	2:D:406:ARG:HH11	2.27	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	133/162 (82%)	123 (92%)	8 (6%)	2 (2%)	13	38
1	B	137/162 (85%)	125 (91%)	11 (8%)	1 (1%)	26	60
1	C	137/162 (85%)	115 (84%)	16 (12%)	6 (4%)	3	9
2	D	9/23 (39%)	9 (100%)	0	0	100	100
2	E	9/23 (39%)	3 (33%)	6 (67%)	0	100	100
2	F	7/23 (30%)	4 (57%)	3 (43%)	0	100	100
All	All	432/555 (78%)	379 (88%)	44 (10%)	9 (2%)	9	28

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	942	ILE
1	C	978	SER
1	A	1013	GLN
1	C	970	GLU
1	C	923	LYS
1	C	941	LYS

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Mol	Chain	Res	Type
1	C	1035	LEU
1	A	1011	SER
1	B	1034	LEU

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	118/142 (83%)	110 (93%)	8 (7%)	20	48
1	B	121/142 (85%)	108 (89%)	13 (11%)	8	23
1	C	121/142 (85%)	103 (85%)	18 (15%)	4	11
2	D	11/22 (50%)	7 (64%)	4 (36%)	0	0
2	E	11/22 (50%)	8 (73%)	3 (27%)	0	1
2	F	9/22 (41%)	7 (78%)	2 (22%)	1	3
All	All	391/492 (80%)	343 (88%)	48 (12%)	6	17

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	940	SER
1	A	978	SER
1	A	988	LYS
1	A	1011	SER
1	A	1013	GLN
1	A	1021	LEU
1	A	1032	LYS
1	A	1049	THR
1	B	921	ASN
1	B	923	LYS
1	B	939	SER
1	B	940	SER
1	B	941	LYS
1	B	948	GLU
1	B	963	THR

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Mol	Chain	Res	Type
1	B	978	SER
1	B	979	THR
1	B	988	LYS
1	B	992	SER
1	B	1021	LEU
1	B	1044	LYS
1	C	927	ASN
1	C	936	ILE
1	C	939	SER
1	C	940	SER
1	C	941	LYS
1	C	942	ILE
1	C	955	LYS
1	C	959	LEU
1	C	962	ARG
1	C	970	GLU
1	C	974	LEU
1	C	979	THR
1	C	988	LYS
1	C	991	ASN
1	C	992	SER
1	C	1032	LYS
1	C	1037	VAL
1	C	1046	LEU
2	D	406	ARG
2	D	409	GLN
2	D	410	ILE
2	D	414	LEU
2	E	406	ARG
2	E	407	MET
2	E	416	GLU
2	F	410	ILE
2	F	416	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	409	GLN

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 carbohydrates 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	134/162 (82%)	-0.46	1 (0%) 89 84	20, 36, 48, 53	0
1	B	139/162 (85%)	-0.42	3 (2%) 65 54	26, 35, 43, 49	0
1	C	139/162 (85%)	-0.26	4 (2%) 55 43	26, 37, 48, 56	0
2	D	11/23 (47%)	2.22	3 (27%) 1 0	75, 78, 83, 83	11 (100%)
2	E	11/23 (47%)	1.46	4 (36%) 0 0	57, 63, 69, 69	11 (100%)
2	F	9/23 (39%)	0.83	1 (11%) 7 3	49, 51, 55, 57	9 (100%)
All	All	443/555 (79%)	-0.24	16 (3%) 46 34	20, 37, 56, 83	31 (6%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	408	SER	8.2
2	E	416	GLU	3.8
2	D	415	SER	3.2
1	B	943	GLN	2.8
1	C	944	PRO	2.7
1	C	948	GLU	2.6
2	E	415	SER	2.5
2	D	409	GLN	2.5
2	E	406	ARG	2.4
1	A	981	ARG	2.3
1	B	1014	GLN	2.3
1	B	940	SER	2.2
1	C	1014	GLN	2.2
2	E	412	ARG	2.1
2	F	415	SER	2.1
1	C	1013	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 carbohydrates 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.