Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 10:21 PM GMT

PDB ID	:	1T62
Title	:	Crystal structure of protein EF3133 from Enterococcus faecalis V583, Pfam DUF984
Authors	:	Fedorov, A.A.; Fedorov, E.V.; Almo, S.C.; Burley, S.K.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on	:	2004-05-05
Resolution	:	3.00 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

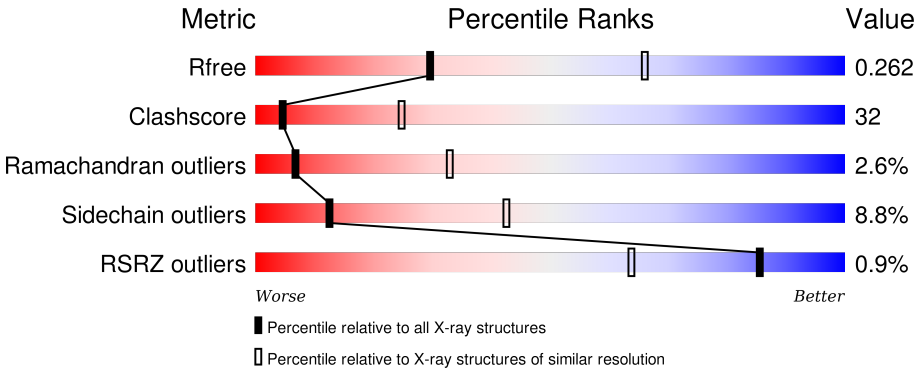
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 3.00 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	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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 <=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	166	% 43% 41% 8% 8%
1	B	166	% 49% 42% 7% •

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2589 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 conserved hypothetical protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	153	Total	C	N	O	S	0	0	0
			1257	805	204	237	11			
1	B	163	Total	C	N	O	S	0	0	0
			1332	849	217	255	11			

There are 20 discrepancies between the modelled and reference sequences:

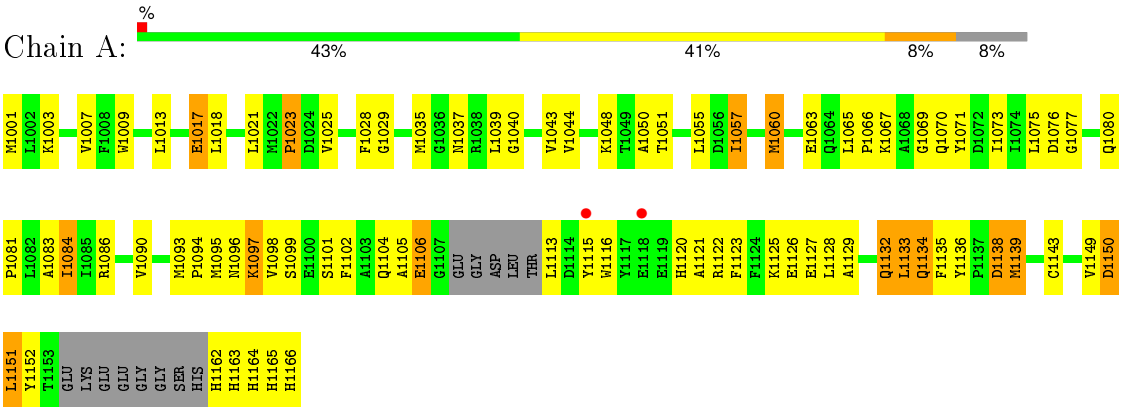
Chain	Residue	Modelled	Actual	Comment	Reference
A	1157	GLU	-	EXPRESSION TAG	UNP Q82ZD1
A	1158	GLY	-	EXPRESSION TAG	UNP Q82ZD1
A	1159	GLY	-	EXPRESSION TAG	UNP Q82ZD1
A	1160	SER	-	EXPRESSION TAG	UNP Q82ZD1
A	1161	HIS	-	EXPRESSION TAG	UNP Q82ZD1
A	1162	HIS	-	EXPRESSION TAG	UNP Q82ZD1
A	1163	HIS	-	EXPRESSION TAG	UNP Q82ZD1
A	1164	HIS	-	EXPRESSION TAG	UNP Q82ZD1
A	1165	HIS	-	EXPRESSION TAG	UNP Q82ZD1
A	1166	HIS	-	EXPRESSION TAG	UNP Q82ZD1
B	2157	GLU	-	EXPRESSION TAG	UNP Q82ZD1
B	2158	GLY	-	EXPRESSION TAG	UNP Q82ZD1
B	2159	GLY	-	EXPRESSION TAG	UNP Q82ZD1
B	2160	SER	-	EXPRESSION TAG	UNP Q82ZD1
B	2161	HIS	-	EXPRESSION TAG	UNP Q82ZD1
B	2162	HIS	-	EXPRESSION TAG	UNP Q82ZD1
B	2163	HIS	-	EXPRESSION TAG	UNP Q82ZD1
B	2164	HIS	-	EXPRESSION TAG	UNP Q82ZD1
B	2165	HIS	-	EXPRESSION TAG	UNP Q82ZD1
B	2166	HIS	-	EXPRESSION TAG	UNP Q82ZD1

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

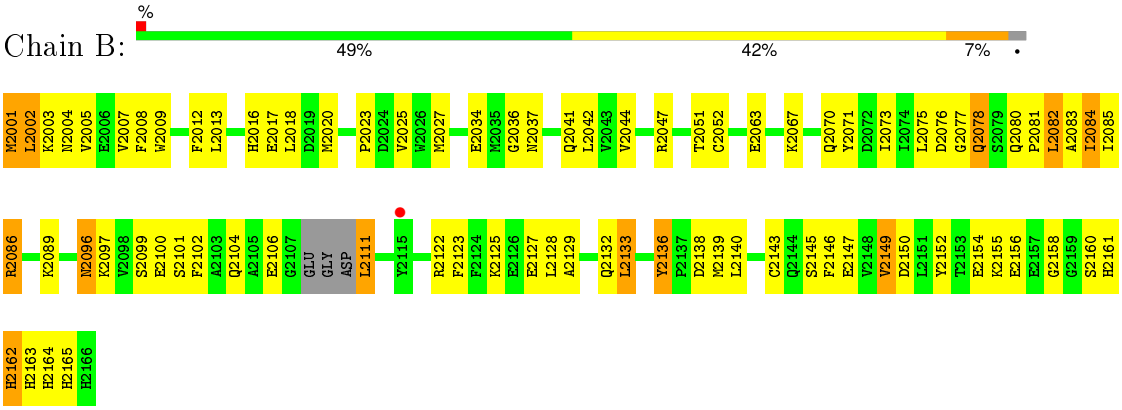
- Molecule 1: conserved hypothetical protein

Chain A:



- Molecule 1: conserved hypothetical protein

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	88.25Å 88.25Å 106.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 3.00 38.21 – 3.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-3.00) 98.8 (38.21-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.89 (at 3.01Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.215 , 0.263 0.212 , 0.262	Depositor DCC
R_{free} test set	520 reflections (5.33%)	DCC
Wilson B-factor (Å^2)	47.4	Xtriage
Anisotropy	0.180	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 75.1	EDS
Estimated twinning fraction	0.035 for -h,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 9889 reflections	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	2589	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.64% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.38	0/1288	0.65	0/1736
1	B	0.43	0/1365	0.64	0/1840
All	All	0.41	0/2653	0.64	0/3576

There are no bond length outliers.

There are no bond angle outliers.

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	1257	0	1199	88	0
1	B	1332	0	1267	87	0
All	All	2589	0	2466	162	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 32.

All (162) 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:B:2160:SER:HA	1:B:2163:HIS:HE1	1.34	0.91
1:B:2037:ASN:ND2	1:B:2077:GLY:H	1.67	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2016:HIS:HB2	1:B:2018:LEU:HD23	1.53	0.90
1:A:1165:HIS:HB3	1:B:2051:THR:HG21	1.54	0.88
1:A:1093:MET:HE1	1:A:1099:SER:H	1.38	0.87
1:A:1037:ASN:ND2	1:A:1077:GLY:H	1.76	0.83
1:A:1025:VAL:HG22	1:B:2001:MET:SD	2.20	0.81
1:B:2084:ILE:HG23	1:B:2150:ASP:HB3	1.66	0.77
1:B:2160:SER:HA	1:B:2163:HIS:CE1	2.19	0.77
1:B:2027:MET:HE3	1:B:2036:GLY:HA3	1.68	0.76
1:A:1029:GLY:HA2	1:A:1035:MET:HE2	1.71	0.72
1:A:1166:HIS:OXT	1:B:2051:THR:HG23	1.90	0.71
1:B:2078:GLN:HG2	1:B:2162:HIS:ND1	2.05	0.71
1:B:2005:VAL:HG11	1:B:2025:VAL:HG11	1.72	0.71
1:A:1129:ALA:HA	1:A:1133:LEU:HD13	1.73	0.70
1:A:1060:MET:CE	1:A:1065:LEU:HD21	2.23	0.69
1:B:2009:TRP:CZ3	1:B:2013:LEU:HD11	2.27	0.69
1:A:1128:LEU:HB3	1:A:1133:LEU:HA	1.74	0.68
1:B:2086:ARG:NH1	1:B:2147:GLU:OE1	2.26	0.68
1:A:1013:LEU:HD22	1:A:1018:LEU:O	1.93	0.68
1:A:1128:LEU:C	1:A:1133:LEU:HB3	2.15	0.66
1:A:1025:VAL:HG12	1:A:1073:ILE:HB	1.78	0.66
1:B:2075:LEU:HD23	1:B:2081:PRO:HA	1.78	0.66
1:A:1066:PRO:O	1:A:1090:VAL:HG21	1.96	0.65
1:A:1069:GLY:HA2	1:A:1086:ARG:HE	1.61	0.65
1:B:2071:TYR:HD2	1:B:2084:ILE:HD11	1.62	0.65
1:A:1037:ASN:ND2	1:A:1076:ASP:HB2	2.12	0.64
1:B:2037:ASN:ND2	1:B:2077:GLY:N	2.44	0.64
1:B:2071:TYR:CD2	1:B:2084:ILE:HD11	2.33	0.63
1:A:1009:TRP:HZ3	1:A:1084:ILE:HD12	1.61	0.63
1:B:2086:ARG:HB3	1:B:2086:ARG:HH11	1.63	0.63
1:A:1113:LEU:HD22	1:A:1115:TYR:HE1	1.63	0.63
1:A:1060:MET:HE1	1:A:1065:LEU:HD21	1.79	0.63
1:A:1003:LYS:O	1:A:1007:VAL:HG23	1.99	0.62
1:A:1163:HIS:HE1	1:B:2063:GLU:OE2	1.82	0.62
1:B:2005:VAL:HG13	1:B:2073:ILE:HG21	1.81	0.62
1:A:1083:ALA:HA	1:A:1151:LEU:HA	1.81	0.62
1:B:2089:LYS:HB3	1:B:2145:SER:HB2	1.80	0.62
1:A:1043:VAL:HG23	1:A:1048:LYS:HD3	1.80	0.61
1:B:2051:THR:HG22	1:B:2052:CYS:H	1.65	0.61
1:A:1163:HIS:CE1	1:B:2063:GLU:OE2	2.54	0.61
1:A:1129:ALA:HA	1:A:1133:LEU:HD22	1.82	0.61
1:B:2128:LEU:HD22	1:B:2133:LEU:HD13	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1165:HIS:HB3	1:B:2051:THR:CG2	2.30	0.60
1:A:1149:VAL:O	1:A:1150:ASP:HB2	2.03	0.59
1:A:1001:MET:HE3	1:A:1080:GLN:HG2	1.83	0.59
1:B:2001:MET:HA	1:B:2004:ASN:OD1	2.03	0.58
1:B:2008:PHE:CE1	1:B:2084:ILE:HG22	2.37	0.58
1:B:2067:LYS:HB3	1:B:2067:LYS:NZ	2.18	0.58
1:A:1037:ASN:HD22	1:A:1077:GLY:H	1.51	0.57
1:B:2084:ILE:HG13	1:B:2149:VAL:HG12	1.87	0.57
1:B:2051:THR:HG22	1:B:2052:CYS:N	2.20	0.56
1:B:2013:LEU:HB3	1:B:2018:LEU:O	2.05	0.56
1:B:2025:VAL:HG12	1:B:2073:ILE:HB	1.88	0.56
1:A:1067:LYS:O	1:A:1070:GLN:HG2	2.06	0.56
1:B:2102:PHE:CE2	1:B:2143:CYS:HB3	2.41	0.56
1:A:1050:ALA:HB3	1:A:1106:GLU:HB3	1.87	0.56
1:B:2076:ASP:HB3	1:B:2082:LEU:HG	1.88	0.56
1:A:1134:GLN:HB2	1:A:1136:TYR:HE1	1.70	0.55
1:B:2009:TRP:HH2	1:B:2020:MET:C	2.10	0.55
1:B:2016:HIS:CB	1:B:2018:LEU:HD23	2.33	0.55
1:B:2016:HIS:HB2	1:B:2018:LEU:CD2	2.34	0.55
1:B:2096:ASN:HD22	1:B:2097:LYS:N	2.04	0.55
1:B:2067:LYS:O	1:B:2070:GLN:HG2	2.06	0.55
1:A:1093:MET:HE2	1:A:1098:VAL:HA	1.90	0.54
1:A:1122:ARG:O	1:A:1126:GLU:HG3	2.07	0.54
1:A:1037:ASN:ND2	1:A:1077:GLY:N	2.50	0.54
1:B:2027:MET:HE3	1:B:2036:GLY:CA	2.37	0.53
1:B:2044:VAL:HG11	1:B:2082:LEU:HD22	1.90	0.53
1:A:1102:PHE:CE2	1:A:1143:CYS:HB3	2.43	0.53
1:B:2096:ASN:HD22	1:B:2096:ASN:C	2.12	0.53
1:A:1127:GLU:HB2	1:B:2164:HIS:HD2	1.73	0.53
1:B:2136:TYR:CE2	1:B:2139:MET:HA	2.43	0.53
1:B:2008:PHE:CZ	1:B:2084:ILE:HG22	2.43	0.53
1:B:2027:MET:CE	1:B:2036:GLY:HA3	2.38	0.53
1:A:1094:PRO:HA	1:A:1139:MET:O	2.09	0.52
1:A:1060:MET:HE3	1:A:1065:LEU:HD21	1.92	0.52
1:B:2005:VAL:HG12	1:B:2073:ILE:HD13	1.92	0.52
1:B:2100:GLU:O	1:B:2104:GLN:HG3	2.11	0.51
1:A:1037:ASN:HD22	1:A:1076:ASP:HB2	1.75	0.50
1:A:1071:TYR:CZ	1:A:1086:ARG:HD2	2.46	0.50
1:A:1121:ALA:O	1:A:1125:LYS:HB2	2.11	0.50
1:A:1001:MET:HG2	1:A:1162:HIS:HE2	1.75	0.50
1:A:1129:ALA:O	1:A:1133:LEU:HD22	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1094:PRO:HD2	1:A:1097:LYS:HD2	1.92	0.50
1:A:1128:LEU:HB3	1:A:1133:LEU:CA	2.42	0.50
1:A:1104:GLN:HG3	1:A:1105:ALA:N	2.25	0.50
1:B:2023:PRO:HB2	1:B:2071:TYR:O	2.12	0.49
1:B:2047:ARG:HG2	1:B:2047:ARG:HH11	1.77	0.49
1:A:1035:MET:HG2	1:A:1039:LEU:HD22	1.93	0.49
1:A:1164:HIS:HB2	1:B:2127:GLU:CD	2.33	0.49
1:B:2002:LEU:HD22	1:B:2002:LEU:O	2.13	0.49
1:A:1063:GLU:OE2	1:B:2160:SER:HB2	2.13	0.49
1:B:2078:GLN:H	1:B:2078:GLN:NE2	2.11	0.49
1:A:1023:PRO:HB2	1:A:1071:TYR:O	2.13	0.48
1:A:1075:LEU:HA	1:A:1080:GLN:O	2.13	0.48
1:A:1083:ALA:HB1	1:A:1150:ASP:O	2.13	0.48
1:B:2136:TYR:CD1	1:B:2136:TYR:N	2.81	0.48
1:B:2158:GLY:HA3	1:B:2161:HIS:HE1	1.79	0.48
1:A:1164:HIS:HB2	1:B:2127:GLU:OE2	2.13	0.48
1:B:2009:TRP:CZ2	1:B:2013:LEU:HD21	2.49	0.47
1:B:2099:SER:OG	1:B:2101:SER:HB3	2.14	0.47
1:A:1040:GLY:O	1:A:1044:VAL:HG23	2.14	0.47
1:A:1057:ILE:HD13	1:A:1057:ILE:O	2.14	0.47
1:A:1151:LEU:N	1:A:1151:LEU:CD2	2.78	0.47
1:A:1093:MET:CE	1:A:1098:VAL:HA	2.44	0.46
1:B:2037:ASN:ND2	1:B:2076:ASP:HB2	2.31	0.46
1:A:1084:ILE:HG23	1:A:1150:ASP:HB3	1.97	0.46
1:A:1028:PHE:HB3	1:A:1039:LEU:HD23	1.98	0.46
1:B:2005:VAL:CG1	1:B:2073:ILE:HG21	2.46	0.46
1:B:2111:LEU:HD23	1:B:2111:LEU:N	2.31	0.46
1:A:1093:MET:HE1	1:A:1099:SER:N	2.17	0.45
1:A:1029:GLY:HA2	1:A:1035:MET:CE	2.43	0.45
1:A:1113:LEU:HD22	1:A:1115:TYR:CE1	2.48	0.45
1:B:2154:GLU:N	1:B:2154:GLU:OE1	2.47	0.45
1:A:1128:LEU:O	1:A:1133:LEU:HB3	2.17	0.45
1:A:1044:VAL:HG13	1:A:1151:LEU:HD12	1.98	0.45
1:A:1055:LEU:HD13	1:A:1055:LEU:C	2.37	0.45
1:B:2003:LYS:O	1:B:2007:VAL:HG23	2.16	0.44
1:B:2078:GLN:O	1:B:2080:GLN:HG2	2.18	0.44
1:A:1075:LEU:HD23	1:A:1081:PRO:HA	2.00	0.44
1:A:1099:SER:C	1:A:1101:SER:N	2.71	0.44
1:A:1009:TRP:CZ2	1:A:1023:PRO:HD3	2.52	0.44
1:B:2102:PHE:HE2	1:B:2143:CYS:HB3	1.81	0.44
1:B:2009:TRP:CH2	1:B:2013:LEU:HD21	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1098:VAL:HG21	1:A:1116:TRP:CH2	2.53	0.43
1:A:1071:TYR:OH	1:A:1086:ARG:NH1	2.48	0.43
1:A:1163:HIS:O	1:A:1164:HIS:HD2	2.02	0.43
1:A:1018:LEU:O	1:A:1018:LEU:HD12	2.18	0.43
1:A:1001:MET:CE	1:A:1080:GLN:HG2	2.47	0.43
1:B:2037:ASN:HD22	1:B:2076:ASP:HB2	1.84	0.43
1:B:2009:TRP:O	1:B:2012:PHE:HB3	2.18	0.43
1:A:1132:GLN:O	1:A:1134:GLN:N	2.45	0.43
1:B:2085:ILE:HD12	1:B:2146:PHE:CG	2.54	0.43
1:A:1096:ASN:HD21	1:A:1138:ASP:HA	1.84	0.42
1:A:1095:MET:HE3	1:A:1120:HIS:HB3	2.01	0.42
1:B:2086:ARG:HB3	1:B:2086:ARG:NH1	2.31	0.42
1:A:1095:MET:HE2	1:A:1095:MET:HB3	1.87	0.42
1:B:2083:ALA:HB1	1:B:2150:ASP:O	2.19	0.42
1:A:1133:LEU:O	1:A:1135:PHE:N	2.53	0.42
1:A:1051:THR:OG1	1:B:2165:HIS:HB3	2.18	0.42
1:A:1081:PRO:HG2	1:A:1152:TYR:CD2	2.54	0.42
1:B:2009:TRP:CH2	1:B:2020:MET:HB2	2.55	0.42
1:A:1093:MET:HA	1:A:1094:PRO:HD3	1.87	0.42
1:B:2078:GLN:HG2	1:B:2162:HIS:HD1	1.79	0.42
1:A:1001:MET:CG	1:A:1162:HIS:HE2	2.33	0.42
1:B:2139:MET:O	1:B:2139:MET:HG3	2.20	0.41
1:B:2152:TYR:CD1	1:B:2155:LYS:HB2	2.55	0.41
1:A:1133:LEU:HD23	1:A:1133:LEU:H	1.85	0.41
1:B:2023:PRO:HG2	1:B:2073:ILE:HG13	2.01	0.41
1:B:2122:ARG:O	1:B:2125:LYS:HB3	2.20	0.41
1:B:2138:ASP:OD2	1:B:2138:ASP:N	2.53	0.41
1:A:1129:ALA:N	1:A:1133:LEU:HB3	2.36	0.41
1:A:1013:LEU:O	1:A:1017:GLU:N	2.53	0.41
1:B:2047:ARG:HG2	1:B:2047:ARG:NH1	2.35	0.41
1:B:2016:HIS:CB	1:B:2018:LEU:CD2	2.96	0.41
1:A:1123:PHE:CZ	1:B:2164:HIS:HB3	2.56	0.41
1:B:2017:GLU:C	1:B:2018:LEU:HD22	2.42	0.40
1:B:2129:ALA:O	1:B:2132:GLN:N	2.54	0.40
1:A:1066:PRO:HA	1:A:1070:GLN:OE1	2.21	0.40
1:B:2096:ASN:ND2	1:B:2096:ASN:C	2.74	0.40
1:A:1164:HIS:HB3	1:B:2123:PHE:CZ	2.57	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	147/166 (89%)	134 (91%)	6 (4%)	7 (5%)	3 17
1	B	159/166 (96%)	150 (94%)	8 (5%)	1 (1%)	30 72
All	All	306/332 (92%)	284 (93%)	14 (5%)	8 (3%)	7 33

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1133	LEU
1	A	1139	MET
1	B	2162	HIS
1	A	1017	GLU
1	A	1060	MET
1	A	1150	ASP
1	A	1097	LYS
1	A	1023	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	138/148 (93%)	130 (94%)	8 (6%)	25 63
1	B	146/148 (99%)	129 (88%)	17 (12%)	7 27
All	All	284/296 (96%)	259 (91%)	25 (9%)	12 42

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1021	LEU
1	A	1057	ILE
1	A	1084	ILE
1	A	1106	GLU
1	A	1132	GLN
1	A	1134	GLN
1	A	1138	ASP
1	A	1151	LEU
1	B	2001	MET
1	B	2002	LEU
1	B	2034	GLU
1	B	2041	GLN
1	B	2042	LEU
1	B	2078	GLN
1	B	2082	LEU
1	B	2084	ILE
1	B	2086	ARG
1	B	2096	ASN
1	B	2106	GLU
1	B	2111	LEU
1	B	2133	LEU
1	B	2136	TYR
1	B	2140	LEU
1	B	2149	VAL
1	B	2156	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1004	ASN
1	A	1037	ASN
1	A	1096	ASN
1	A	1163	HIS
1	B	2037	ASN
1	B	2041	GLN
1	B	2078	GLN
1	B	2080	GLN
1	B	2096	ASN
1	B	2134	GLN
1	B	2163	HIS
1	B	2164	HIS

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	153/166 (92%)	-0.24	2 (1%)	79	53	10, 37, 76, 106	0
1	B	163/166 (98%)	-0.51	1 (0%)	90	73	8, 26, 60, 96	0
All	All	316/332 (95%)	-0.38	3 (0%)	85	64	8, 30, 69, 106	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1115	TYR	2.8
1	A	1118	GLU	2.4
1	B	2115	TYR	2.4

6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.