



Full wwPDB X-ray Structure Validation Report

May 13, 2020 – 01:31 am BST

PDB ID : 4GML
Title : Crystal structure of human NOT1 MIF4G domain
Authors : Petit, P.; Weichenrieder, O.; Wohlbold, L.; Izaurralde, E.
Deposited on : 2012-08-16
Resolution : 2.90 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 following versions of software and data (see [references](#) ) were used in the production of this report:

MolProbity : 4.02b-467
Xtrriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

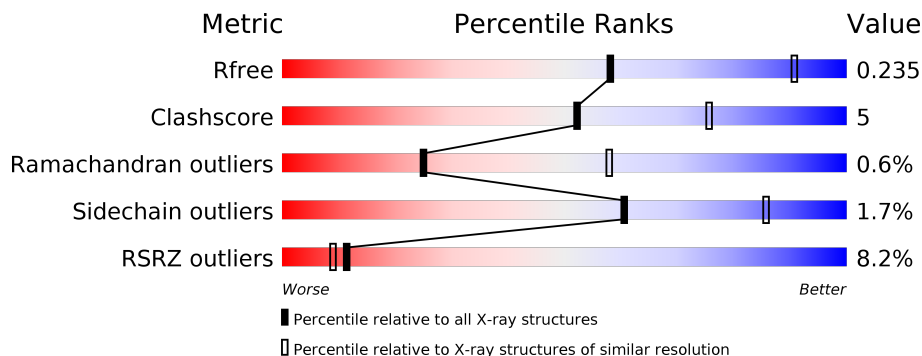
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 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	235	 3% 90% 9%
1	B	235	 9% 80% 18%
1	C	235	 9% 80% 13% 5%
1	D	235	 3% 88% 11%
1	E	235	 14% 84% 12%
1	F	235	 11% 80% 18%

2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 11132 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 CCR4-NOT transcription complex subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	235	Total 1896	C 1224	N 315	O 348	S 9	0	0	0
1	B	235	Total 1893	C 1222	N 315	O 347	S 9	0	0	0
1	C	223	Total 1786	C 1154	N 293	O 331	S 8	0	0	0
1	D	233	Total 1871	C 1207	N 312	O 343	S 9	0	0	0
1	E	226	Total 1811	C 1166	N 301	O 336	S 8	0	0	0
1	F	233	Total 1845	C 1190	N 306	O 340	S 9	0	0	0

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	VAL	-	EXPRESSION TAG	UNP A5YKK6
A	-8	LEU	-	EXPRESSION TAG	UNP A5YKK6
A	-7	PHE	-	EXPRESSION TAG	UNP A5YKK6
A	-6	GLN	-	EXPRESSION TAG	UNP A5YKK6
A	-5	GLY	-	EXPRESSION TAG	UNP A5YKK6
A	-4	PRO	-	EXPRESSION TAG	UNP A5YKK6
A	-3	HIS	-	EXPRESSION TAG	UNP A5YKK6
A	-2	MET	-	EXPRESSION TAG	UNP A5YKK6
A	-1	LEU	-	EXPRESSION TAG	UNP A5YKK6
A	0	GLU	-	EXPRESSION TAG	UNP A5YKK6
B	-9	VAL	-	EXPRESSION TAG	UNP A5YKK6
B	-8	LEU	-	EXPRESSION TAG	UNP A5YKK6
B	-7	PHE	-	EXPRESSION TAG	UNP A5YKK6
B	-6	GLN	-	EXPRESSION TAG	UNP A5YKK6
B	-5	GLY	-	EXPRESSION TAG	UNP A5YKK6
B	-4	PRO	-	EXPRESSION TAG	UNP A5YKK6
B	-3	HIS	-	EXPRESSION TAG	UNP A5YKK6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	MET	-	EXPRESSION TAG	UNP A5YKK6
B	-1	LEU	-	EXPRESSION TAG	UNP A5YKK6
B	0	GLU	-	EXPRESSION TAG	UNP A5YKK6
C	-9	VAL	-	EXPRESSION TAG	UNP A5YKK6
C	-8	LEU	-	EXPRESSION TAG	UNP A5YKK6
C	-7	PHE	-	EXPRESSION TAG	UNP A5YKK6
C	-6	GLN	-	EXPRESSION TAG	UNP A5YKK6
C	-5	GLY	-	EXPRESSION TAG	UNP A5YKK6
C	-4	PRO	-	EXPRESSION TAG	UNP A5YKK6
C	-3	HIS	-	EXPRESSION TAG	UNP A5YKK6
C	-2	MET	-	EXPRESSION TAG	UNP A5YKK6
C	-1	LEU	-	EXPRESSION TAG	UNP A5YKK6
C	0	GLU	-	EXPRESSION TAG	UNP A5YKK6
D	-9	VAL	-	EXPRESSION TAG	UNP A5YKK6
D	-8	LEU	-	EXPRESSION TAG	UNP A5YKK6
D	-7	PHE	-	EXPRESSION TAG	UNP A5YKK6
D	-6	GLN	-	EXPRESSION TAG	UNP A5YKK6
D	-5	GLY	-	EXPRESSION TAG	UNP A5YKK6
D	-4	PRO	-	EXPRESSION TAG	UNP A5YKK6
D	-3	HIS	-	EXPRESSION TAG	UNP A5YKK6
D	-2	MET	-	EXPRESSION TAG	UNP A5YKK6
D	-1	LEU	-	EXPRESSION TAG	UNP A5YKK6
D	0	GLU	-	EXPRESSION TAG	UNP A5YKK6
E	-9	VAL	-	EXPRESSION TAG	UNP A5YKK6
E	-8	LEU	-	EXPRESSION TAG	UNP A5YKK6
E	-7	PHE	-	EXPRESSION TAG	UNP A5YKK6
E	-6	GLN	-	EXPRESSION TAG	UNP A5YKK6
E	-5	GLY	-	EXPRESSION TAG	UNP A5YKK6
E	-4	PRO	-	EXPRESSION TAG	UNP A5YKK6
E	-3	HIS	-	EXPRESSION TAG	UNP A5YKK6
E	-2	MET	-	EXPRESSION TAG	UNP A5YKK6
E	-1	LEU	-	EXPRESSION TAG	UNP A5YKK6
E	0	GLU	-	EXPRESSION TAG	UNP A5YKK6
F	-9	VAL	-	EXPRESSION TAG	UNP A5YKK6
F	-8	LEU	-	EXPRESSION TAG	UNP A5YKK6
F	-7	PHE	-	EXPRESSION TAG	UNP A5YKK6
F	-6	GLN	-	EXPRESSION TAG	UNP A5YKK6
F	-5	GLY	-	EXPRESSION TAG	UNP A5YKK6
F	-4	PRO	-	EXPRESSION TAG	UNP A5YKK6
F	-3	HIS	-	EXPRESSION TAG	UNP A5YKK6
F	-2	MET	-	EXPRESSION TAG	UNP A5YKK6
F	-1	LEU	-	EXPRESSION TAG	UNP A5YKK6

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Chain	Residue	Modelled	Actual	Comment	Reference
F	0	GLU	-	EXPRESSION TAG	UNP A5YKK6

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	9	Total O 9 9	0	0
2	B	6	Total O 6 6	0	0
2	C	4	Total O 4 4	0	0
2	D	8	Total O 8 8	0	0
2	E	1	Total O 1 1	0	0
2	F	2	Total O 2 2	0	0

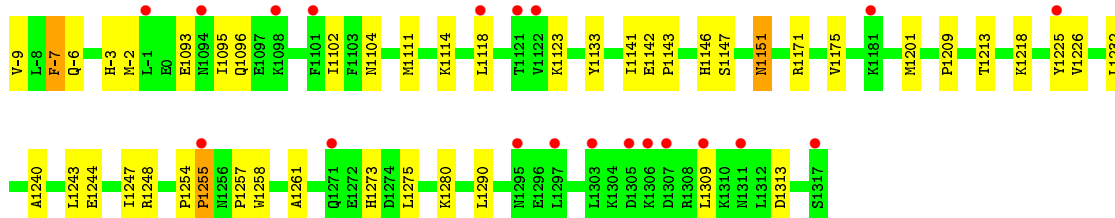
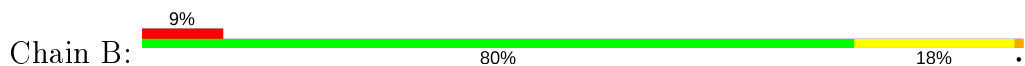
3 Residue-property plots [i](#)

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

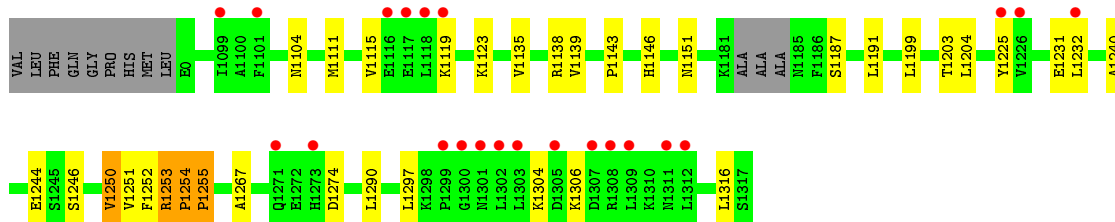
- Molecule 1: CCR4-NOT transcription complex subunit 1



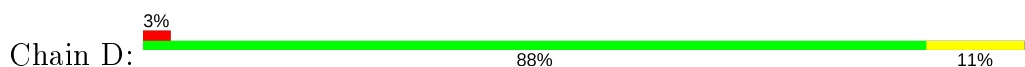
- Molecule 1: CCR4-NOT transcription complex subunit 1



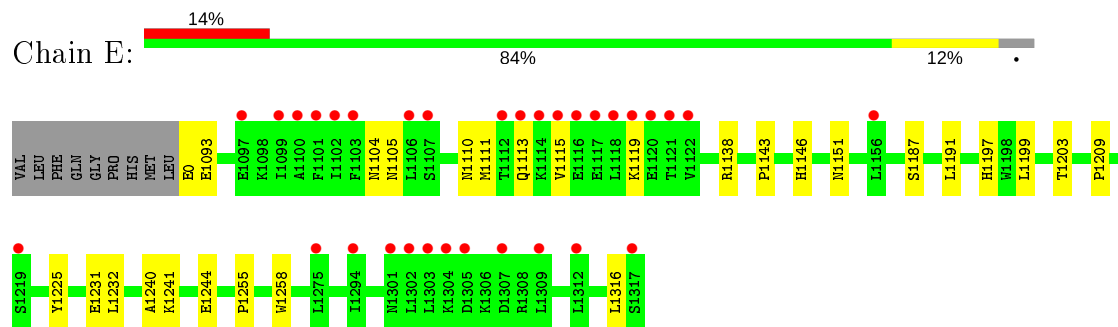
- Molecule 1: CCR4-NOT transcription complex subunit 1



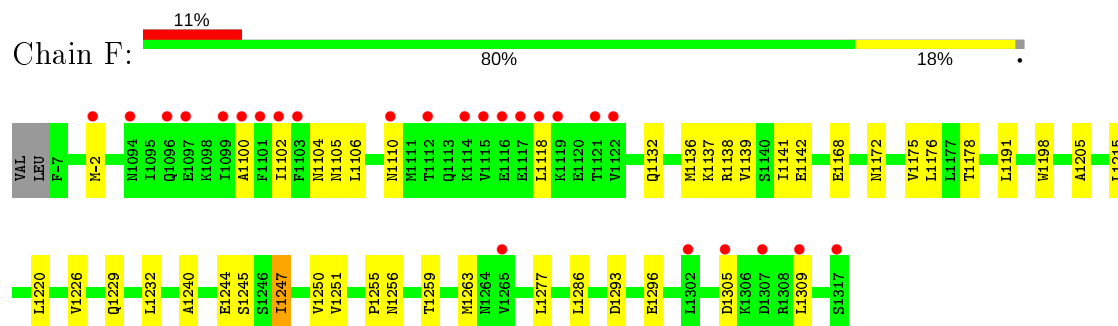
- Molecule 1: CCR4-NOT transcription complex subunit 1



● Molecule 1: CCR4-NOT transcription complex subunit 1



● Molecule 1: CCR4-NOT transcription complex subunit 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	88.27Å 133.82Å 331.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.31 – 2.90 49.31 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.31-2.90) 95.2 (49.31-2.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 2.91Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.190 , 0.229 0.195 , 0.235	Depositor DCC
R_{free} test set	2216 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	61.8	Xtrriage
Anisotropy	0.765	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 85.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11132	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.23	0/1933	0.40	0/2618
1	B	0.23	0/1930	0.40	0/2615
1	C	0.24	0/1819	0.39	0/2465
1	D	0.23	0/1907	0.38	0/2582
1	E	0.23	0/1845	0.41	0/2501
1	F	0.23	0/1881	0.40	0/2555
All	All	0.23	0/11315	0.40	0/15336

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 [i](#)

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	1896	0	1930	14	0
1	B	1893	0	1923	31	0
1	C	1786	0	1797	25	0
1	D	1871	0	1898	19	0
1	E	1811	0	1825	16	0
1	F	1845	0	1840	33	0
2	A	9	0	0	0	0
2	B	6	0	0	0	0
2	C	4	0	0	0	0
2	D	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	1	0	0	0	0
2	F	2	0	0	0	0
All	All	11132	0	11213	122	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 5.

All (122) 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:1257:PRO:HG2	1:C:1253:ARG:HH21	1.32	0.92
1:D:1203:THR:HG23	1:D:1204:LEU:N	1.82	0.92
1:F:1100:ALA:O	1:F:1104:ASN:HB2	1.73	0.88
1:D:1203:THR:HG23	1:D:1204:LEU:H	1.39	0.87
1:C:1246:SER:HB2	1:C:1252:PHE:O	1.76	0.85
1:F:1139:VAL:HG11	1:F:1198:TRP:CG	2.13	0.83
1:D:1203:THR:CG2	1:D:1204:LEU:H	1.95	0.79
1:D:1203:THR:CG2	1:D:1204:LEU:N	2.48	0.76
1:D:1104:ASN:O	1:D:1114:LYS:NZ	2.20	0.75
1:B:-3:HIS:HB3	1:F:1191:LEU:HD13	1.71	0.73
1:A:1094:ASN:HB3	1:A:1097:GLU:HB3	1.71	0.72
1:B:1104:ASN:O	1:B:1114:LYS:NZ	2.25	0.69
1:D:1203:THR:HG23	1:D:1204:LEU:HG	1.75	0.68
1:C:1225:TYR:HA	1:C:1232:LEU:HD11	1.76	0.68
1:D:1200:GLY:HA2	1:D:1203:THR:HG22	1.76	0.68
1:F:1102:ILE:HG13	1:F:1118:LEU:HD23	1.77	0.66
1:F:1100:ALA:O	1:F:1104:ASN:CB	2.44	0.64
1:F:1141:ILE:HG13	1:F:1142:GLU:HG3	1.79	0.64
1:B:-9:VAL:HG21	1:B:1133:TYR:HB2	1.80	0.63
1:F:1104:ASN:O	1:F:1105:ASN:HB2	1.98	0.63
1:C:1111:MET:SD	1:C:1151:ASN:ND2	2.73	0.61
1:B:1225:TYR:HA	1:B:1232:LEU:HD11	1.83	0.61
1:B:1258:TRP:HB2	1:C:1250:VAL:HG21	1.83	0.60
1:D:1260:MET:HE1	1:D:1292:LEU:HD11	1.84	0.59
1:A:-9:VAL:HG21	1:A:1133:TYR:HB2	1.84	0.59
1:C:1199:LEU:O	1:C:1203:THR:HG22	2.01	0.59
1:B:1141:ILE:HG13	1:B:1142:GLU:HG3	1.84	0.59
1:B:1123:LYS:HE3	1:F:1178:THR:HG21	1.84	0.59
1:B:1257:PRO:HG2	1:C:1253:ARG:NH2	2.12	0.58
1:E:1258:TRP:HB2	1:F:1250:VAL:HG11	1.86	0.57
1:B:1247:ILE:HD11	1:B:1290:LEU:HA	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1104:ASN:OD1	1:C:1138:ARG:NH2	2.36	0.57
1:D:1202:ILE:HA	1:D:1206:LYS:HE3	1.87	0.56
1:E:1104:ASN:OD1	1:E:1138:ARG:NH2	2.34	0.56
1:A:1260:MET:HE1	1:A:1292:LEU:HD11	1.87	0.56
1:E:1209:PRO:HB3	1:F:1251:VAL:HG21	1.86	0.56
1:F:1215:LEU:HD11	1:F:1220:LEU:HD22	1.87	0.56
1:F:1100:ALA:O	1:F:1104:ASN:N	2.37	0.55
1:F:1205:ALA:HB2	1:F:1245:SER:HB3	1.87	0.55
1:E:1111:MET:HE1	1:E:1151:ASN:C	2.27	0.54
1:F:1106:LEU:HD12	1:F:1110:ASN:OD1	2.07	0.54
1:F:1136:MET:HG2	1:F:1172:ASN:OD1	2.09	0.53
1:C:1203:THR:HG23	1:C:1204:LEU:H	1.74	0.53
1:A:1138:ARG:NH2	1:A:1142:GLU:OE2	2.42	0.53
1:C:1231:GLU:HG2	1:C:1316:LEU:HD13	1.89	0.53
1:F:1251:VAL:O	1:F:1256:ASN:ND2	2.39	0.52
1:B:1111:MET:SD	1:B:1151:ASN:HB3	2.50	0.52
1:B:-2:MET:HE1	1:F:1175:VAL:HG11	1.90	0.52
1:D:1135:VAL:HA	1:D:1139:VAL:HB	1.91	0.52
1:E:1209:PRO:HD3	1:F:1251:VAL:HG11	1.91	0.51
1:E:1231:GLU:HG2	1:E:1316:LEU:HD13	1.93	0.51
1:A:1183:ALA:O	1:A:1189:ARG:NH2	2.44	0.51
1:E:1225:TYR:HA	1:E:1232:LEU:HD21	1.92	0.51
1:B:1209:PRO:HB3	1:C:1251:VAL:HG21	1.93	0.50
1:B:1093:GLU:HB3	1:B:1095:ILE:HG13	1.93	0.50
1:C:1253:ARG:O	1:C:1254:PRO:C	2.50	0.50
1:D:1200:GLY:CA	1:D:1203:THR:HG22	2.42	0.50
1:C:1115:VAL:O	1:C:1119:LYS:HG2	2.11	0.50
1:B:1226:VAL:HG11	1:B:1309:LEU:HB3	1.93	0.49
1:D:1114:LYS:HA	1:D:1117:GLU:HG2	1.94	0.49
1:B:1275:LEU:HD23	1:B:1280:LYS:HG2	1.95	0.48
1:B:1275:LEU:O	1:B:1280:LYS:NZ	2.34	0.48
1:B:-6:GLN:O	1:F:1137:LYS:NZ	2.46	0.48
1:F:1259:THR:O	1:F:1263:MET:HG2	2.13	0.48
1:A:1144:ASN:OD1	1:B:1248:ARG:NH2	2.42	0.48
1:D:1200:GLY:HA2	1:D:1203:THR:CG2	2.43	0.47
1:B:-3:HIS:HB2	1:F:1176:LEU:HD21	1.96	0.47
1:E:1143:PRO:HA	1:E:1146:HIS:CD2	2.49	0.47
1:A:-6:GLN:NE2	1:A:-2:MET:O	2.48	0.47
1:E:1187:SER:O	1:E:1191:LEU:HG	2.15	0.47
1:D:1102:ILE:O	1:D:1114:LYS:HD3	2.15	0.46
1:B:1143:PRO:HA	1:B:1146:HIS:CE1	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1267:ALA:HB1	1:D:1299:PRO:HG3	1.98	0.46
1:C:1240:ALA:O	1:C:1244:GLU:HG3	2.16	0.46
1:A:1251:VAL:O	1:A:1256:ASN:ND2	2.33	0.45
1:F:1293:ASP:HB3	1:F:1296:GLU:HB2	1.98	0.45
1:C:1135:VAL:HA	1:C:1139:VAL:HB	1.98	0.45
1:F:1229:GLN:HA	1:F:1232:LEU:HD12	1.99	0.45
1:A:1133:TYR:CZ	1:A:1138:ARG:HD3	2.52	0.45
1:B:1254:PRO:HA	1:B:1255:PRO:HA	1.81	0.45
1:D:1305:ASP:C	1:D:1307:ASP:H	2.20	0.45
1:C:1203:THR:HG23	1:C:1204:LEU:HG	1.99	0.44
1:E:1197:HIS:HB2	1:E:1241:LYS:HD2	1.99	0.44
1:F:1240:ALA:O	1:F:1244:GLU:HG3	2.18	0.44
1:A:1240:ALA:O	1:A:1244:GLU:HG3	2.17	0.44
1:B:1102:ILE:HG21	1:B:1118:LEU:HB2	2.00	0.44
1:C:1123:LYS:HB3	1:C:1123:LYS:HE2	1.76	0.44
1:C:1254:PRO:HA	1:C:1255:PRO:HA	1.81	0.44
1:F:1139:VAL:CG1	1:F:1198:TRP:HB2	2.47	0.44
1:F:1247:ILE:HD13	1:F:1247:ILE:HA	1.75	0.44
1:E:1115:VAL:O	1:E:1119:LYS:HG3	2.17	0.44
1:A:1181:LYS:NZ	1:A:1317:SER:OG	2.51	0.44
1:E:1240:ALA:O	1:E:1244:GLU:HG3	2.18	0.43
1:C:1253:ARG:O	1:C:1254:PRO:O	2.36	0.43
1:E:1258:TRP:CB	1:F:1250:VAL:HG11	2.49	0.43
1:C:1203:THR:HG23	1:C:1204:LEU:N	2.34	0.43
1:C:1253:ARG:HH11	1:C:1253:ARG:HB2	1.84	0.43
1:D:1240:ALA:O	1:D:1244:GLU:HG3	2.19	0.42
1:D:1254:PRO:HA	1:D:1255:PRO:HA	1.79	0.42
1:B:1218:LYS:HE3	1:B:1261:ALA:HB1	2.01	0.42
1:B:1240:ALA:O	1:B:1244:GLU:HG3	2.19	0.42
1:B:1147:SER:HA	1:B:1201:MET:HE1	2.02	0.42
1:D:1226:VAL:HG11	1:D:1309:LEU:HB3	2.00	0.42
1:E:1241:LYS:NZ	1:E:1244:GLU:OE2	2.50	0.42
1:B:1243:LEU:HD23	1:B:1243:LEU:HA	1.87	0.42
1:B:-7:PHE:O	1:F:1136:MET:HE2	2.19	0.42
1:B:-9:VAL:HG13	1:B:1096:GLN:HG3	2.01	0.42
1:F:1226:VAL:HG11	1:F:1309:LEU:HG	2.02	0.42
1:A:1093:GLU:O	1:A:1093:GLU:HG3	2.19	0.42
1:F:1305:ASP:O	1:F:1309:LEU:HB2	2.21	0.41
1:F:1138:ARG:HA	1:F:1138:ARG:HD2	1.96	0.41
1:A:-2:MET:HE2	1:A:-2:MET:HB2	1.92	0.41
1:B:-2:MET:HG3	1:F:1136:MET:HE3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:1132:GLN:HA	1:F:1168:GLU:HG2	2.02	0.41
1:C:1143:PRO:HA	1:C:1146:HIS:CD2	2.56	0.41
1:A:-8:LEU:HB2	1:A:1132:GLN:OE1	2.21	0.41
1:B:1171:ARG:O	1:B:1175:VAL:HG23	2.20	0.41
1:C:1304:LYS:O	1:C:1306:LYS:N	2.51	0.41
1:E:1110:ASN:O	1:E:1113:GLN:HG2	2.21	0.41
1:E:1199:LEU:O	1:E:1203:THR:OG1	2.31	0.40
1:C:1187:SER:O	1:C:1191:LEU:HG	2.21	0.40
1:C:1267:ALA:HB2	1:C:1297:LEU:HD13	2.03	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	233/235 (99%)	226 (97%)	7 (3%)	0	100	100
1	B	233/235 (99%)	224 (96%)	7 (3%)	2 (1%)	17	48
1	C	219/235 (93%)	208 (95%)	9 (4%)	2 (1%)	17	48
1	D	231/235 (98%)	221 (96%)	9 (4%)	1 (0%)	34	66
1	E	224/235 (95%)	214 (96%)	8 (4%)	2 (1%)	17	48
1	F	231/235 (98%)	222 (96%)	8 (4%)	1 (0%)	34	66
All	All	1371/1410 (97%)	1315 (96%)	48 (4%)	8 (1%)	25	58

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	1254	PRO
1	B	-7	PHE
1	E	1105	ASN

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Mol	Chain	Res	Type
1	D	1255	PRO
1	C	1255	PRO
1	B	1255	PRO
1	E	1255	PRO
1	F	1255	PRO

5.3.2 Protein sidechains [i](#)

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	215/220 (98%)	211 (98%)	4 (2%)	57 84
1	B	214/220 (97%)	210 (98%)	4 (2%)	57 84
1	C	201/220 (91%)	197 (98%)	4 (2%)	55 82
1	D	210/220 (96%)	207 (99%)	3 (1%)	67 89
1	E	203/220 (92%)	201 (99%)	2 (1%)	76 92
1	F	204/220 (93%)	200 (98%)	4 (2%)	55 82
All	All	1247/1320 (94%)	1226 (98%)	21 (2%)	60 86

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-8	LEU
1	A	1093	GLU
1	A	1094	ASN
1	A	1250	VAL
1	B	1151	ASN
1	B	1213	THR
1	B	1273	HIS
1	B	1313	ASP
1	C	1250	VAL
1	C	1253	ARG
1	C	1274	ASP
1	C	1290	LEU
1	D	-7	PHE

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Mol	Chain	Res	Type
1	D	1151	ASN
1	D	1253	ARG
1	E	0	GLU
1	E	1093	GLU
1	F	-2	MET
1	F	1247	ILE
1	F	1277	LEU
1	F	1286	LEU

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

Mol	Chain	Res	Type
1	F	1311	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 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	235/235 (100%)	0.32	8 (3%) 45 40	44, 66, 104, 123	0
1	B	235/235 (100%)	0.48	20 (8%) 10 8	49, 71, 112, 142	0
1	C	223/235 (94%)	0.48	22 (9%) 7 5	24, 69, 113, 129	0
1	D	233/235 (99%)	0.37	7 (3%) 50 45	28, 65, 108, 137	0
1	E	226/235 (96%)	0.64	32 (14%) 2 2	55, 79, 119, 152	0
1	F	233/235 (99%)	0.57	25 (10%) 6 4	54, 77, 117, 130	0
All	All	1385/1410 (98%)	0.47	114 (8%) 11 9	24, 72, 115, 152	0

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	1102	ILE	5.2
1	E	1275	LEU	4.8
1	E	1117	GLU	4.8
1	E	1101	PHE	4.6
1	F	1121	THR	4.6
1	E	1116	GLU	4.2
1	C	1309	LEU	4.2
1	C	1226	VAL	4.0
1	F	1102	ILE	4.0
1	E	1114	LYS	3.9
1	A	1116	GLU	3.9
1	B	1271	GLN	3.8
1	E	1120	GLU	3.7
1	E	1113	GLN	3.7
1	E	1317	SER	3.6
1	C	1305	ASP	3.6
1	F	-2	MET	3.5
1	B	1309	LEU	3.5
1	E	1106	LEU	3.5

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Mol	Chain	Res	Type	RSRZ
1	C	1301	ASN	3.5
1	F	1317	SER	3.5
1	E	1309	LEU	3.4
1	F	1103	PHE	3.4
1	E	1121	THR	3.4
1	B	1225	TYR	3.4
1	F	1117	GLU	3.3
1	A	-9	VAL	3.3
1	E	1294	ILE	3.3
1	B	-1	LEU	3.3
1	F	1114	LYS	3.3
1	E	1118	LEU	3.3
1	F	1101	PHE	3.3
1	C	1271	GLN	3.2
1	E	1103	PHE	3.2
1	E	1304	LYS	3.1
1	E	1312	LEU	3.1
1	F	1100	ALA	3.1
1	F	1096	GLN	3.1
1	B	1307	ASP	3.1
1	C	1303	LEU	3.1
1	C	1117	GLU	3.0
1	E	1100	ALA	3.0
1	A	1307	ASP	3.0
1	B	1317	SER	3.0
1	F	1118	LEU	3.0
1	D	1303	LEU	2.9
1	F	1112	THR	2.9
1	C	1312	LEU	2.9
1	E	1301	ASN	2.8
1	E	1107	SER	2.8
1	E	1302	LEU	2.8
1	F	1305	ASP	2.8
1	F	1307	ASP	2.8
1	B	1122	VAL	2.7
1	B	1303	LEU	2.7
1	F	1099	ILE	2.7
1	B	1297	LEU	2.7
1	E	1122	VAL	2.7
1	E	1305	ASP	2.7
1	C	1307	ASP	2.6
1	D	1309	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	1311	ASN	2.5
1	B	1101	PHE	2.5
1	C	1116	GLU	2.5
1	F	1097	GLU	2.5
1	F	1122	VAL	2.5
1	C	1099	ILE	2.5
1	D	1310	LYS	2.5
1	F	1094	ASN	2.5
1	D	1182	ALA	2.4
1	F	1115	VAL	2.4
1	B	1094	ASN	2.4
1	C	1118	LEU	2.4
1	A	1317	SER	2.4
1	A	1271	GLN	2.4
1	C	1225	TYR	2.4
1	F	1110	ASN	2.4
1	E	1099	ILE	2.3
1	B	1306	LYS	2.3
1	C	1308	ARG	2.3
1	E	1115	VAL	2.3
1	F	1302	LEU	2.3
1	E	1112	THR	2.3
1	C	1273	HIS	2.3
1	F	1116	GLU	2.3
1	C	1299	PRO	2.3
1	C	1119	LYS	2.3
1	E	1307	ASP	2.2
1	A	1101	PHE	2.2
1	C	1232	LEU	2.2
1	E	1119	LYS	2.2
1	B	1098	LYS	2.2
1	B	1118	LEU	2.2
1	C	1302	LEU	2.2
1	D	1101	PHE	2.2
1	B	1255	PRO	2.2
1	B	1181	LYS	2.2
1	C	1101	PHE	2.1
1	A	1309	LEU	2.1
1	B	1305	ASP	2.1
1	F	1309	LEU	2.1
1	B	1295	ASN	2.1
1	B	1121	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	F	1265	VAL	2.1
1	A	1118	LEU	2.1
1	E	1097	GLU	2.1
1	E	1303	LEU	2.1
1	D	1311	ASN	2.0
1	E	1156	LEU	2.0
1	E	1219	SER	2.0
1	C	1300	GLY	2.0
1	F	1119	LYS	2.0
1	D	1295	ASN	2.0
1	C	1311	ASN	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.