Full wwPDB X-ray Structure Validation Report (i)

Jun 17, 2024 – 02:50 AM EDT

PDB ID	:	3UF1
Title	:	Crystal Structure of Vimentin (fragment 144-251) from Homo sapiens, Northeast Structural Genomics Consortium Target HR4796B
Authors	:	Kuzin, A.; Abashidze, M.; Vorobiev, S.M.; Patel, P.; Xiao, R.; Ciccosanti, C.; Shastry, R.; Everett, J.K.; Nair, R.; Acton, T.B.; Rost, B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on	:	2011-10-31
Resolution	:	2.81 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 (i) 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 (i)) were used in the production of this report:

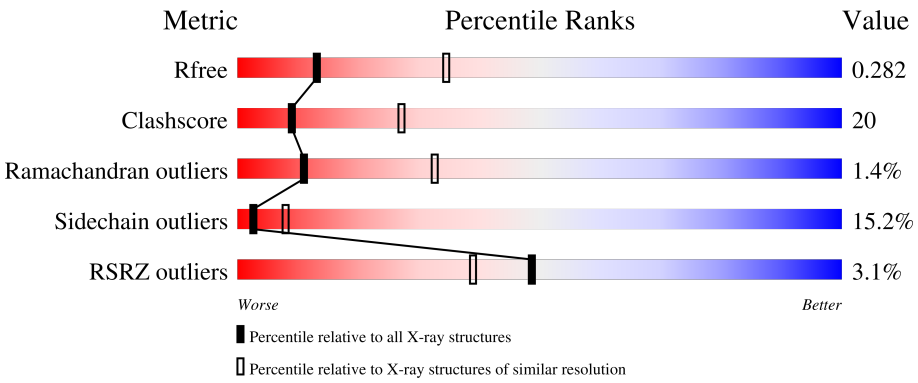
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.81 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	3617 (2.84-2.80)
Clashscore	141614	4060 (2.84-2.80)
Ramachandran outliers	138981	3978 (2.84-2.80)
Sidechain outliers	138945	3980 (2.84-2.80)
RSRZ outliers	127900	3552 (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 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 <=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	119	2%, 50%, 32%, 6%, 13%
1	B	119	3%, 59%, 24%, 6%, 11%
1	C	119	3%, 52%, 34%, 6%, 8%
1	D	119	3%, 61%, 26%, 5%, •, 6%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 3639 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 Vimentin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	104	875	532	158	182	3	0	0	0
1	B	106	892	541	163	185	3	0	0	0
1	C	109	924	559	173	188	4	0	0	0
1	D	112	948	573	179	192	4	0	0	0

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	133	MSE	-	EXPRESSION TAG	UNP P08670
A	134	GLY	-	EXPRESSION TAG	UNP P08670
A	135	HIS	-	EXPRESSION TAG	UNP P08670
A	136	HIS	-	EXPRESSION TAG	UNP P08670
A	137	HIS	-	EXPRESSION TAG	UNP P08670
A	138	HIS	-	EXPRESSION TAG	UNP P08670
A	139	HIS	-	EXPRESSION TAG	UNP P08670
A	140	HIS	-	EXPRESSION TAG	UNP P08670
A	141	SER	-	EXPRESSION TAG	UNP P08670
A	142	HIS	-	EXPRESSION TAG	UNP P08670
A	143	MSE	-	EXPRESSION TAG	UNP P08670
B	133	MSE	-	EXPRESSION TAG	UNP P08670
B	134	GLY	-	EXPRESSION TAG	UNP P08670
B	135	HIS	-	EXPRESSION TAG	UNP P08670
B	136	HIS	-	EXPRESSION TAG	UNP P08670
B	137	HIS	-	EXPRESSION TAG	UNP P08670
B	138	HIS	-	EXPRESSION TAG	UNP P08670
B	139	HIS	-	EXPRESSION TAG	UNP P08670
B	140	HIS	-	EXPRESSION TAG	UNP P08670
B	141	SER	-	EXPRESSION TAG	UNP P08670
B	142	HIS	-	EXPRESSION TAG	UNP P08670

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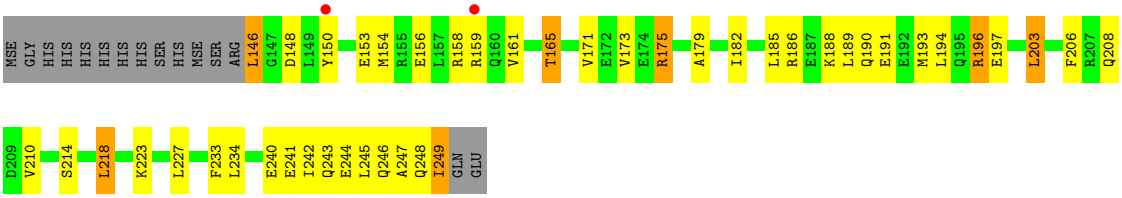
Chain	Residue	Modelled	Actual	Comment	Reference
B	143	MSE	-	EXPRESSION TAG	UNP P08670
C	133	MSE	-	EXPRESSION TAG	UNP P08670
C	134	GLY	-	EXPRESSION TAG	UNP P08670
C	135	HIS	-	EXPRESSION TAG	UNP P08670
C	136	HIS	-	EXPRESSION TAG	UNP P08670
C	137	HIS	-	EXPRESSION TAG	UNP P08670
C	138	HIS	-	EXPRESSION TAG	UNP P08670
C	139	HIS	-	EXPRESSION TAG	UNP P08670
C	140	HIS	-	EXPRESSION TAG	UNP P08670
C	141	SER	-	EXPRESSION TAG	UNP P08670
C	142	HIS	-	EXPRESSION TAG	UNP P08670
C	143	MSE	-	EXPRESSION TAG	UNP P08670
D	133	MSE	-	EXPRESSION TAG	UNP P08670
D	134	GLY	-	EXPRESSION TAG	UNP P08670
D	135	HIS	-	EXPRESSION TAG	UNP P08670
D	136	HIS	-	EXPRESSION TAG	UNP P08670
D	137	HIS	-	EXPRESSION TAG	UNP P08670
D	138	HIS	-	EXPRESSION TAG	UNP P08670
D	139	HIS	-	EXPRESSION TAG	UNP P08670
D	140	HIS	-	EXPRESSION TAG	UNP P08670
D	141	SER	-	EXPRESSION TAG	UNP P08670
D	142	HIS	-	EXPRESSION TAG	UNP P08670
D	143	MSE	-	EXPRESSION TAG	UNP P08670

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.

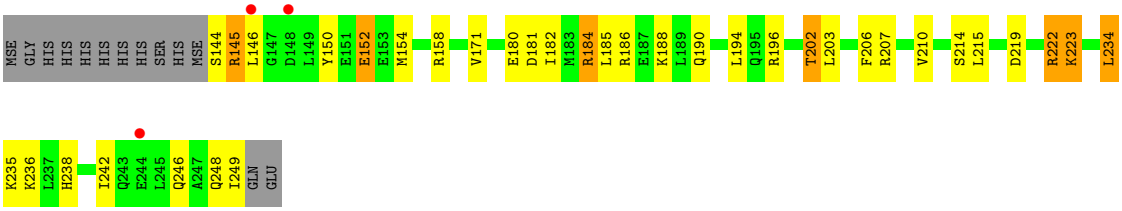
• Molecule 1: Vimentin

Chain A:



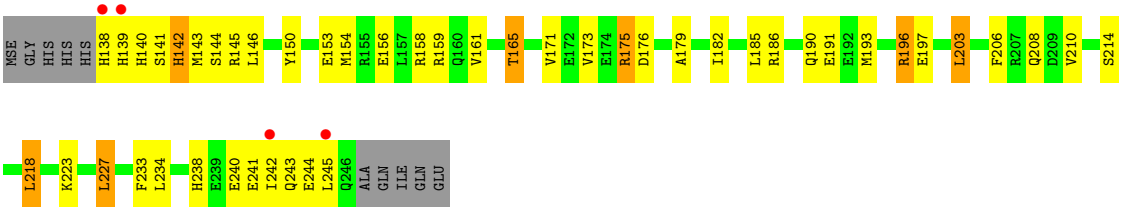
• Molecule 1: Vimentin

Chain B:



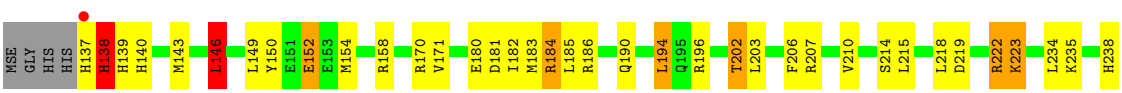
• Molecule 1: Vimentin

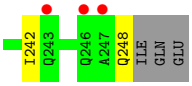
Chain C:



• Molecule 1: Vimentin

Chain D:




I242
Q243
Q246
A247
Q248
ILE
GLN
GLU

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	61.02Å 86.81Å 114.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.80 – 2.81 29.49 – 2.81	Depositor EDS
% Data completeness (in resolution range)	98.7 (19.80-2.81) 99.0 (29.49-2.81)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	6.96 (at 2.80Å)	Xtriage
Refinement program	PHENIX 1.7.1_743	Depositor
R, R_{free}	0.234 , 0.284 0.233 , 0.282	Depositor DCC
R_{free} test set	1500 reflections (9.82%)	wwPDB-VP
Wilson B-factor (Å^2)	53.2	Xtriage
Anisotropy	0.897	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 57.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	3639	wwPDB-VP
Average B, all atoms (Å^2)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.19% 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, 0.333 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	0.43	0/875	0.59	0/1164
1	B	0.46	0/892	0.55	0/1186
1	C	0.41	0/927	0.58	0/1231
1	D	0.40	0/952	0.55	1/1265 (0.1%)
All	All	0.42	0/3646	0.57	1/4846 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	146	LEU	CA-CB-CG	5.23	127.33	115.30

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	875	0	862	46	0
1	B	892	0	880	39	0
1	C	924	0	898	53	0
1	D	948	0	918	48	0
All	All	3639	0	3558	145	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 20.

All (145) 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:175:ARG:HH11	1:C:175:ARG:HG2	1.14	1.12
1:B:154:MSE:HE3	1:B:158:ARG:HH22	1.13	1.11
1:A:175:ARG:HH11	1:A:175:ARG:HG2	1.14	1.05
1:C:143:MSE:HE1	1:D:143:MSE:HG3	1.40	1.03
1:D:154:MSE:HE3	1:D:158:ARG:HH22	1.29	0.93
1:A:171:VAL:HG12	1:B:171:VAL:HG12	1.56	0.86
1:C:171:VAL:HG12	1:D:171:VAL:HG12	1.59	0.84
1:C:159:ARG:HH11	1:C:159:ARG:HG2	1.45	0.82
1:B:207:ARG:HG3	1:B:207:ARG:HH11	1.45	0.81
1:D:207:ARG:HH11	1:D:207:ARG:HG3	1.46	0.81
1:B:154:MSE:HE3	1:B:158:ARG:NH2	1.94	0.79
1:A:159:ARG:HG2	1:A:159:ARG:HH11	1.47	0.78
1:A:175:ARG:HG2	1:A:175:ARG:NH1	1.95	0.78
1:C:142:HIS:ND1	1:C:143:MSE:HA	2.02	0.75
1:C:175:ARG:HH11	1:C:175:ARG:CG	1.99	0.74
1:A:246:GLN:O	1:A:249:ILE:HG12	1.89	0.73
1:B:154:MSE:CE	1:B:158:ARG:HH22	1.98	0.73
1:A:175:ARG:HH11	1:A:175:ARG:CG	1.97	0.72
1:C:210:VAL:HG22	1:D:210:VAL:HG23	1.71	0.70
1:D:181:ASP:HA	1:D:184:ARG:HH11	1.56	0.70
1:A:210:VAL:HG22	1:B:210:VAL:HG23	1.74	0.69
1:C:138:HIS:N	1:C:139:HIS:HA	2.07	0.69
1:D:137:HIS:CG	1:D:138:HIS:N	2.62	0.68
1:C:242:ILE:HG13	1:D:238:HIS:HE1	1.57	0.68
1:C:143:MSE:HE3	1:D:139:HIS:CD2	2.30	0.67
1:C:206:PHE:CE2	1:D:206:PHE:HB3	2.31	0.66
1:C:206:PHE:HE2	1:D:206:PHE:HB3	1.62	0.65
1:A:171:VAL:HG12	1:B:171:VAL:CG1	2.24	0.65
1:D:137:HIS:O	1:D:139:HIS:N	2.29	0.65
1:B:181:ASP:HA	1:B:184:ARG:HH11	1.60	0.64
1:A:203:LEU:HD11	1:B:202:THR:HG21	1.80	0.64
1:D:181:ASP:OD1	1:D:184:ARG:NH1	2.31	0.64
1:C:175:ARG:HG2	1:C:175:ARG:NH1	1.95	0.61
1:B:246:GLN:O	1:B:249:ILE:HG23	2.01	0.61
1:C:234:LEU:HD11	1:D:235:LYS:HE3	1.83	0.61
1:B:207:ARG:HG3	1:B:207:ARG:NH1	2.16	0.61
1:C:142:HIS:HA	1:C:144:SER:H	1.67	0.60
1:A:242:ILE:HG13	1:B:238:HIS:HE1	1.66	0.60
1:C:146:LEU:HD23	1:D:146:LEU:HD11	1.84	0.60
1:A:247:ALA:O	1:A:248:GLN:HB2	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:203:LEU:O	1:C:203:LEU:HD22	2.03	0.59
1:A:233:PHE:CE2	1:D:154:MSE:HE1	2.37	0.59
1:A:234:LEU:HD11	1:B:235:LYS:HE3	1.86	0.58
1:C:142:HIS:CG	1:C:143:MSE:HA	2.39	0.58
1:C:171:VAL:HG12	1:D:171:VAL:CG1	2.30	0.58
1:D:223:LYS:O	1:D:223:LYS:HG3	2.05	0.57
1:A:206:PHE:CE2	1:B:206:PHE:HB3	2.39	0.57
1:A:206:PHE:HE2	1:B:206:PHE:HB3	1.70	0.56
1:A:203:LEU:HD22	1:A:203:LEU:O	2.06	0.56
1:A:240:GLU:OE1	1:C:145:ARG:HB3	2.06	0.55
1:B:223:LYS:O	1:B:223:LYS:HG3	2.06	0.55
1:C:185:LEU:HD21	1:D:186:ARG:HG2	1.88	0.55
1:C:150:TYR:CG	1:D:150:TYR:CD2	2.95	0.55
1:C:203:LEU:HD11	1:D:202:THR:HG21	1.88	0.54
1:C:242:ILE:HG13	1:D:238:HIS:CE1	2.42	0.54
1:A:179:ALA:O	1:A:182:ILE:HG22	2.08	0.54
1:D:137:HIS:CE1	1:D:138:HIS:HB2	2.42	0.53
1:A:233:PHE:CD2	1:D:154:MSE:HE1	2.44	0.53
1:A:182:ILE:HG13	1:A:186:ARG:NH2	2.23	0.53
1:A:182:ILE:HD13	1:B:182:ILE:HG13	1.90	0.53
1:C:182:ILE:HD13	1:D:182:ILE:HG13	1.91	0.53
1:B:180:GLU:HG3	1:B:184:ARG:HH12	1.73	0.52
1:B:238:HIS:O	1:B:242:ILE:HG23	2.09	0.52
1:C:142:HIS:CB	1:C:143:MSE:HA	2.40	0.52
1:A:214:SER:O	1:A:218:LEU:HD22	2.10	0.52
1:A:189:LEU:HD21	1:B:188:LYS:HD3	1.91	0.51
1:C:159:ARG:HH11	1:C:159:ARG:CG	2.19	0.51
1:D:137:HIS:CG	1:D:138:HIS:H	2.27	0.51
1:C:179:ALA:O	1:C:182:ILE:HG22	2.10	0.51
1:D:181:ASP:HA	1:D:184:ARG:NH1	2.25	0.51
1:A:154:MSE:HE3	1:A:158:ARG:HH21	1.76	0.51
1:A:159:ARG:HH11	1:A:159:ARG:CG	2.21	0.51
1:D:207:ARG:HG3	1:D:207:ARG:NH1	2.17	0.50
1:C:182:ILE:HG13	1:C:186:ARG:NH2	2.26	0.50
1:A:161:VAL:O	1:A:165:THR:HB	2.11	0.50
1:C:161:VAL:O	1:C:165:THR:HB	2.12	0.49
1:B:181:ASP:OD1	1:B:184:ARG:NH1	2.45	0.49
1:C:138:HIS:N	1:C:139:HIS:CA	2.75	0.49
1:C:244:GLU:O	1:C:245:LEU:HD23	2.11	0.49
1:D:238:HIS:O	1:D:242:ILE:HG23	2.12	0.49
1:C:214:SER:O	1:C:218:LEU:HD22	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:LEU:HD21	1:B:186:ARG:HG2	1.95	0.49
1:A:150:TYR:O	1:A:154:MSE:HG3	2.13	0.49
1:A:185:LEU:HD23	1:B:185:LEU:HB2	1.95	0.47
1:C:234:LEU:HD23	1:C:234:LEU:HA	1.70	0.47
1:A:244:GLU:O	1:A:245:LEU:HD23	2.14	0.47
1:C:185:LEU:HD23	1:D:185:LEU:HB2	1.96	0.47
1:C:240:GLU:C	1:C:242:ILE:H	2.18	0.47
1:D:180:GLU:HG3	1:D:184:ARG:HH12	1.79	0.47
1:D:190:GLN:HA	1:D:190:GLN:OE1	2.15	0.47
1:C:154:MSE:HE3	1:C:158:ARG:HH21	1.80	0.47
1:C:175:ARG:CG	1:C:175:ARG:NH1	2.67	0.47
1:D:154:MSE:CE	1:D:158:ARG:HH22	2.14	0.47
1:B:207:ARG:HH11	1:B:207:ARG:CG	2.23	0.47
1:B:181:ASP:HA	1:B:184:ARG:NH1	2.30	0.46
1:A:242:ILE:HG13	1:B:238:HIS:CE1	2.48	0.46
1:C:150:TYR:O	1:C:154:MSE:HG3	2.16	0.46
1:A:185:LEU:HD23	1:B:185:LEU:CB	2.45	0.45
1:A:240:GLU:OE1	1:C:145:ARG:NE	2.49	0.45
1:B:145:ARG:NE	1:B:145:ARG:HA	2.32	0.45
1:C:223:LYS:NZ	1:C:223:LYS:HB2	2.31	0.45
1:C:150:TYR:CG	1:D:150:TYR:HD2	2.34	0.45
1:C:159:ARG:HG2	1:C:159:ARG:NH1	2.20	0.45
1:D:222:ARG:HD3	1:D:222:ARG:HA	1.75	0.45
1:B:207:ARG:NH1	1:B:207:ARG:CG	2.81	0.44
1:B:152:GLU:HA	1:B:152:GLU:OE1	2.17	0.44
1:B:203:LEU:C	1:B:203:LEU:HD13	2.37	0.44
1:D:154:MSE:HE3	1:D:158:ARG:NH2	2.13	0.44
1:A:240:GLU:C	1:A:242:ILE:H	2.21	0.44
1:B:154:MSE:HE1	1:C:233:PHE:CE2	2.53	0.44
1:C:175:ARG:NH1	1:C:176:ASP:OD2	2.48	0.44
1:D:170:ARG:NH1	1:D:170:ARG:HG3	2.33	0.44
1:D:152:GLU:HA	1:D:152:GLU:OE1	2.17	0.44
1:A:182:ILE:HD13	1:B:182:ILE:CG1	2.48	0.43
1:B:222:ARG:HD3	1:B:222:ARG:HA	1.70	0.43
1:D:137:HIS:O	1:D:138:HIS:C	2.57	0.43
1:A:159:ARG:HG2	1:A:159:ARG:NH1	2.21	0.43
1:C:185:LEU:HD23	1:D:185:LEU:CB	2.48	0.43
1:A:175:ARG:NH1	1:A:175:ARG:CG	2.66	0.43
1:A:223:LYS:HB2	1:A:223:LYS:NZ	2.33	0.43
1:C:196:ARG:HH21	1:C:197:GLU:HB2	1.83	0.43
1:A:247:ALA:O	1:A:248:GLN:CB	2.64	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:190:GLN:OE1	1:B:190:GLN:HA	2.19	0.43
1:C:240:GLU:O	1:C:242:ILE:N	2.51	0.43
1:C:153:GLU:HA	1:C:153:GLU:OE1	2.18	0.42
1:B:154:MSE:HE1	1:C:233:PHE:CD2	2.55	0.42
1:D:203:LEU:C	1:D:203:LEU:HD13	2.40	0.42
1:D:207:ARG:NH1	1:D:207:ARG:CG	2.81	0.42
1:C:143:MSE:H	1:C:143:MSE:HG3	1.71	0.42
1:A:150:TYR:CG	1:B:150:TYR:CD2	3.08	0.41
1:D:146:LEU:HA	1:D:149:LEU:HB2	2.00	0.41
1:A:153:GLU:HA	1:A:153:GLU:OE1	2.20	0.41
1:C:159:ARG:CG	1:C:159:ARG:NH1	2.79	0.41
1:D:170:ARG:HG3	1:D:170:ARG:HH11	1.85	0.41
1:A:146:LEU:C	1:A:148:ASP:N	2.73	0.41
1:B:234:LEU:HA	1:B:234:LEU:HD22	1.82	0.41
1:A:159:ARG:CG	1:A:159:ARG:NH1	2.81	0.41
1:A:194:LEU:HD23	1:A:194:LEU:HA	1.84	0.41
1:A:196:ARG:HH21	1:A:197:GLU:HB2	1.86	0.41
1:D:143:MSE:O	1:D:146:LEU:HD12	2.21	0.41
1:D:194:LEU:HD22	1:D:194:LEU:HA	1.87	0.41
1:C:227:LEU:HD12	1:C:227:LEU:HA	1.85	0.40
1:A:171:VAL:CG1	1:B:171:VAL:HG12	2.39	0.40
1:D:140:HIS:CD2	1:D:140:HIS:O	2.74	0.40
1:D:218:LEU:HD23	1:D:218:LEU:HA	1.93	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	102/119 (86%)	93 (91%)	7 (7%)	2 (2%)	7 23
1	B	104/119 (87%)	99 (95%)	5 (5%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	C	107/119 (90%)	97 (91%)	7 (6%)	3 (3%)	5 15
1	D	110/119 (92%)	107 (97%)	2 (2%)	1 (1%)	17 44
All	All	423/476 (89%)	396 (94%)	21 (5%)	6 (1%)	11 32

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	141	SER
1	D	138	HIS
1	C	140	HIS
1	A	241	GLU
1	C	241	GLU
1	A	188	LYS

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	96/105 (91%)	81 (84%)	15 (16%)	2 7
1	B	98/105 (93%)	82 (84%)	16 (16%)	2 7
1	C	102/105 (97%)	87 (85%)	15 (15%)	3 9
1	D	104/105 (99%)	89 (86%)	15 (14%)	3 9
All	All	400/420 (95%)	339 (85%)	61 (15%)	2 8

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	146	LEU
1	A	156	GLU
1	A	165	THR
1	A	173	VAL
1	A	175	ARG
1	A	190	GLN

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Mol	Chain	Res	Type
1	A	191	GLU
1	A	193	MSE
1	A	196	ARG
1	A	203	LEU
1	A	208	GLN
1	A	218	LEU
1	A	227	LEU
1	A	243	GLN
1	A	249	ILE
1	B	144	SER
1	B	145	ARG
1	B	146	LEU
1	B	152	GLU
1	B	184	ARG
1	B	194	LEU
1	B	196	ARG
1	B	202	THR
1	B	214	SER
1	B	215	LEU
1	B	219	ASP
1	B	222	ARG
1	B	223	LYS
1	B	234	LEU
1	B	236	LYS
1	B	248	GLN
1	C	142	HIS
1	C	156	GLU
1	C	165	THR
1	C	173	VAL
1	C	175	ARG
1	C	190	GLN
1	C	191	GLU
1	C	193	MSE
1	C	196	ARG
1	C	203	LEU
1	C	208	GLN
1	C	218	LEU
1	C	227	LEU
1	C	238	HIS
1	C	243	GLN
1	D	138	HIS
1	D	146	LEU

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Mol	Chain	Res	Type
1	D	152	GLU
1	D	183	MSE
1	D	184	ARG
1	D	194	LEU
1	D	196	ARG
1	D	202	THR
1	D	214	SER
1	D	215	LEU
1	D	219	ASP
1	D	222	ARG
1	D	223	LYS
1	D	234	LEU
1	D	248	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	139	HIS
1	D	238	HIS

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 monosaccharides 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 (i)

6.1 Protein, DNA and RNA chains (i)

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	101/119 (84%)	-0.10	2 (1%)	65	56	39, 67, 133, 157	0
1	B	103/119 (86%)	-0.08	3 (2%)	51	41	36, 63, 152, 194	0
1	C	105/119 (88%)	0.03	4 (3%)	40	30	43, 76, 143, 183	0
1	D	108/119 (90%)	0.10	4 (3%)	41	31	44, 72, 168, 221	1 (0%)
All	All	417/476 (87%)	-0.01	13 (3%)	49	39	36, 71, 145, 221	1 (0%)

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	148	ASP	4.8
1	C	138	HIS	4.5
1	D	246	GLN	4.0
1	D	247	ALA	3.7
1	D	137	HIS	3.5
1	C	242	ILE	3.4
1	B	146	LEU	3.2
1	C	245	LEU	3.1
1	D	243	GLN	2.8
1	A	150	TYR	2.7
1	C	139	HIS	2.5
1	A	159	ARG	2.3
1	B	244	GLU	2.2

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 monosaccharides 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.