



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 5, 2021 – 10:02 AM EDT

PDB ID : 7N94  
Title : LINE-1 endonuclease domain complex with DNA  
Authors : Korolev, S.; Miller, I.  
Deposited on : 2021-06-16  
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](mailto: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.

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

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.23.2  
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) : 2.23.2

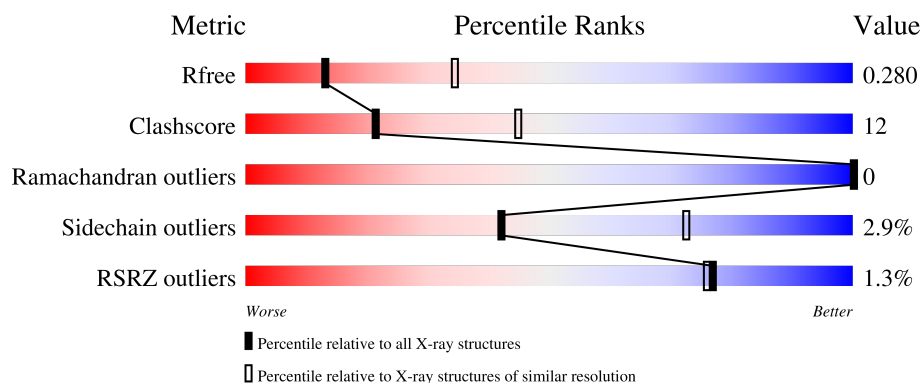
# 1 Overall quality at a glance

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))
$R_{free}$	130704	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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  $\geq 3$ , 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	238	<div> <div>2%</div> <div>67%</div> <div>29%</div> <div>..</div> </div>
1	B	238	<div> <div>73%</div> <div>23%</div> <div>..</div> </div>
2	D	18	<div> <div>11%</div> <div>78%</div> <div>22%</div> </div>
2	F	18	<div> <div>83%</div> <div>11%</div> <div>6%</div> </div>
3	E	18	<div> <div>6%</div> <div>67%</div> <div>28%</div> <div>6%</div> </div>

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Mol	Chain	Length	Quality of chain
3	G	18	 78% 22%

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5149 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 LINE-1 retrotransposable element ORF2 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	232	Total	C	N	O	S	0	0	0
			1860	1171	342	341	6			
1	B	232	Total	C	N	O	S	0	0	0
			1860	1171	342	341	6			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	15	ILE	VAL	conflict	UNP O00370
A	21	ALA	PRO	conflict	UNP O00370
A	145	ALA	ASP	engineered mutation	UNP O00370
A	152	THR	ILE	conflict	UNP O00370
A	175	ALA	THR	conflict	UNP O00370
A	226	LYS	TYR	engineered mutation	UNP O00370
B	15	ILE	VAL	conflict	UNP O00370
B	21	ALA	PRO	conflict	UNP O00370
B	145	ALA	ASP	engineered mutation	UNP O00370
B	152	THR	ILE	conflict	UNP O00370
B	175	ALA	THR	conflict	UNP O00370
B	226	LYS	TYR	engineered mutation	UNP O00370

- Molecule 2 is a DNA chain called DNA (5'-D(\*GP\*CP\*TP\*CP\*CP\*TP\*TP\*TP\*TP\*AP\*AP\*GP\*GP\*GP\*CP\*TP\*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	18	Total	C	N	O	P	0	0	0
			364	176	61	110	17			
2	F	17	Total	C	N	O	P	0	0	0
			343	166	56	105	16			

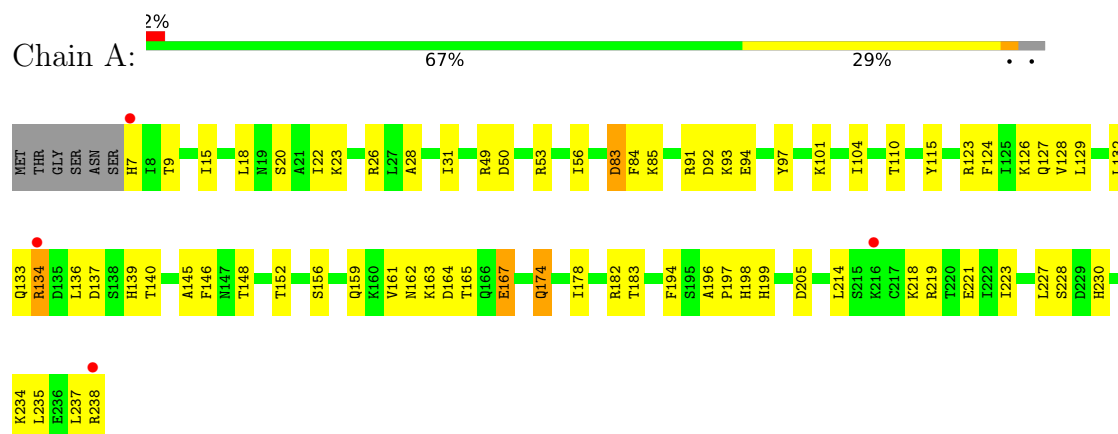
- Molecule 3 is a DNA chain called DNA (5'-D(\*AP\*GP\*CP\*CP\*CP\*TP\*TP\*AP\*AP\*AP\*AP\*AP\*GP\*GP\*AP\*GP\*CP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	17	Total 351	C 166	N 71	O 97	P 17	0	0	0
3	G	18	Total 371	C 176	N 73	O 104	P 18	0	0	0

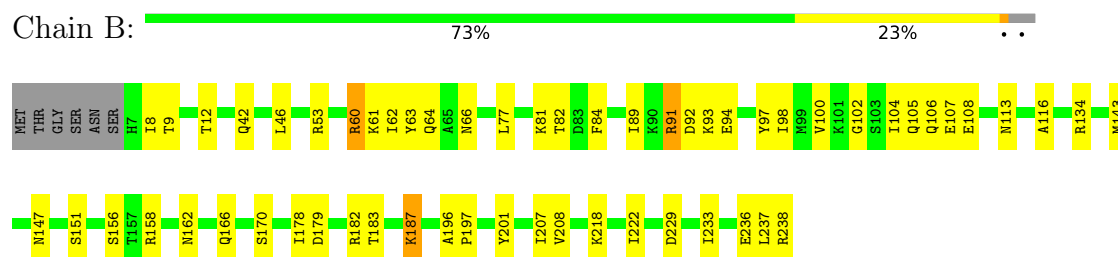
### 3 Residue-property plots

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: LINE-1 retrotransposable element ORF2 protein

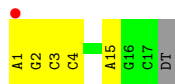


- Molecule 1: LINE-1 retrotransposable element ORF2 protein

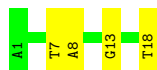
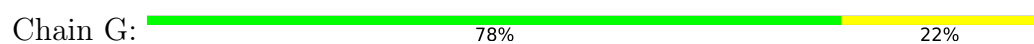




- Molecule 3: DNA (5'-D(\*AP\*GP\*CP\*CP\*CP\*TP\*TP\*AP\*AP\*AP\*AP\*AP\*GP\*GP\*AP\*GP\*CP\*T)-3')



- Molecule 3: DNA (5'-D(\*AP\*GP\*CP\*CP\*CP\*TP\*TP\*AP\*AP\*AP\*AP\*AP\*GP\*GP\*AP\*GP\*CP\*T)-3')



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	100.77 Å   100.77 Å   140.77 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	42.92 – 2.85 42.92 – 2.85	Depositor EDS
% Data completeness (in resolution range)	99.3 (42.92-2.85) 96.8 (42.92-2.85)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.04 (at 2.86 Å)	Xtriage
Refinement program	PHENIX 1.16_3549, PHENIX 1.16_3549	Depositor
R, $R_{free}$	0.214 , 0.280 0.214 , 0.280	Depositor DCC
$R_{free}$ test set	1744 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	43.5	Xtriage
Anisotropy	0.023	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 35.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	5149	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.42	0/1893	0.76	3/2552 (0.1%)
1	B	0.41	0/1893	0.74	4/2552 (0.2%)
2	D	0.67	0/406	0.97	0/625
2	F	0.83	0/382	1.04	0/588
3	E	0.71	0/395	0.96	0/607
3	G	0.84	0/417	0.96	0/641
All	All	0.54	0/5386	0.83	7/7565 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	60	ARG	CG-CD-NE	-10.21	90.35	111.80
1	B	81	LYS	CD-CE-NZ	-8.85	91.36	111.70
1	A	174	GLN	CA-CB-CG	6.94	128.67	113.40
1	B	187	LYS	CD-CE-NZ	6.49	126.63	111.70
1	A	167	GLU	CA-CB-CG	6.42	127.52	113.40
1	A	83	ASP	CB-CG-OD1	6.07	123.77	118.30
1	B	60	ARG	NE-CZ-NH2	-5.38	117.61	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	1860	0	1924	68	1
1	B	1860	0	1924	38	0
2	D	364	0	207	4	0
2	F	343	0	196	1	0
3	E	351	0	190	5	0
3	G	371	0	202	4	0
All	All	5149	0	4643	117	1

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 12.

All (117) 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:A:7:HIS:CD2	1:A:238:ARG:HA	1.66	1.30
1:A:7:HIS:HD2	1:A:238:ARG:HA	0.90	1.05
1:A:7:HIS:HD2	1:A:238:ARG:CA	1.70	1.04
1:A:133:GLN:HA	1:A:136:LEU:HD12	1.54	0.89
1:A:7:HIS:CD2	1:A:238:ARG:CA	2.52	0.85
1:B:222:ILE:HG12	1:B:233:ILE:HD13	1.62	0.80
1:B:9:THR:HG22	1:B:236:GLU:HG2	1.65	0.78
1:B:92:ASP:OD2	1:B:94:GLU:N	2.15	0.78
1:A:7:HIS:HB3	1:A:237:LEU:O	1.86	0.75
1:B:89:ILE:HG12	1:B:100:VAL:HG12	1.69	0.74
1:B:64:GLN:O	1:B:91:ARG:NH1	2.21	0.74
1:A:7:HIS:CD2	1:A:238:ARG:HG2	2.25	0.71
1:A:133:GLN:CD	1:A:134:ARG:HH21	1.94	0.71
1:A:84:PHE:HD1	1:A:104:ILE:HG13	1.56	0.71
1:B:8:ILE:HG23	1:B:237:LEU:HB2	1.76	0.68
1:A:164:ASP:HA	1:A:167:GLU:HB2	1.76	0.68
1:B:62:ILE:HG12	1:B:77:LEU:HD22	1.76	0.67
1:A:235:LEU:HG	1:A:237:LEU:HD13	1.76	0.67
3:G:18:DT:H6	3:G:18:DT:OP2	1.78	0.67
1:A:49:ARG:HD3	3:G:13:DG:OP2	1.95	0.66
2:D:17:DT:H72	3:E:2:DG:N2	2.11	0.66
1:B:53:ARG:HH11	1:B:53:ARG:HG2	1.61	0.66
1:A:126:LYS:HD2	1:A:167:GLU:O	1.98	0.64
1:A:92:ASP:OD2	1:A:94:GLU:N	2.22	0.63
1:B:12:THR:HG23	1:B:42:GLN:HE22	1.63	0.63
1:A:92:ASP:OD2	1:A:93:LYS:N	2.32	0.62
1:B:116:ALA:O	1:B:147:ASN:ND2	2.31	0.62
1:B:143:MET:HG2	1:B:207:ILE:HG23	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:ASP:CG	1:A:139:HIS:H	2.04	0.60
3:E:3:DC:H2''	3:E:4:DC:O5'	2.01	0.60
1:B:92:ASP:HB2	1:B:97:TYR:CE2	2.37	0.60
1:B:106:GLN:O	1:B:107:GLU:HG2	2.02	0.59
1:A:197:PRO:HB2	1:A:198:HIS:HD2	1.66	0.59
1:A:9:THR:HG21	1:A:234:LYS:HE3	1.85	0.58
1:B:218:LYS:HE2	1:B:238:ARG:HD3	1.85	0.58
1:B:91:ARG:HB3	1:B:98:ILE:HD12	1.86	0.57
1:A:164:ASP:O	1:A:167:GLU:HB3	2.05	0.57
1:A:161:VAL:HG12	1:A:165:THR:HB	1.87	0.56
1:B:92:ASP:OD2	1:B:93:LYS:N	2.38	0.56
1:A:93:LYS:HA	1:A:93:LYS:HE2	1.87	0.56
1:A:178:ILE:HD11	1:A:214:LEU:HD13	1.88	0.55
1:A:133:GLN:NE2	1:A:134:ARG:HH21	2.05	0.55
1:B:92:ASP:OD1	1:B:94:GLU:HB2	2.08	0.54
1:B:207:ILE:HD11	1:B:233:ILE:HG22	1.90	0.53
1:B:222:ILE:HG12	1:B:233:ILE:CD1	2.37	0.53
1:A:137:ASP:OD1	1:A:140:THR:N	2.41	0.53
1:A:7:HIS:CD2	1:A:238:ARG:CB	2.92	0.52
1:A:7:HIS:NE2	1:A:238:ARG:HG2	2.24	0.52
1:A:137:ASP:CG	1:A:140:THR:HG23	2.31	0.52
1:B:104:ILE:HG22	1:B:105:GLN:HG3	1.92	0.52
1:A:31:ILE:HG21	1:A:56:ILE:HG21	1.93	0.51
1:A:123:ARG:HD2	1:A:164:ASP:OD2	2.10	0.51
1:A:197:PRO:HB2	1:A:198:HIS:CD2	2.46	0.50
1:B:82:THR:HG23	1:B:104:ILE:HG23	1.94	0.50
2:D:2:DC:H2''	2:D:3:DT:H5'	1.93	0.49
1:A:84:PHE:CD1	1:A:104:ILE:HG13	2.43	0.49
1:A:235:LEU:HG	1:A:237:LEU:CD1	2.42	0.49
3:E:3:DC:H2'	3:E:4:DC:C5	2.48	0.49
1:A:84:PHE:HD1	1:A:104:ILE:CG1	2.25	0.49
1:A:219:ARG:NH2	1:A:221:GLU:OE1	2.45	0.49
1:B:53:ARG:HG2	1:B:53:ARG:NH1	2.28	0.49
1:A:137:ASP:OD1	1:A:140:THR:HG23	2.13	0.48
3:G:18:DT:OP2	3:G:18:DT:C6	2.64	0.48
1:A:28:ALA:HA	1:A:56:ILE:HD12	1.96	0.47
1:A:164:ASP:HA	1:A:167:GLU:CB	2.44	0.47
1:B:179:ASP:HA	1:B:208:VAL:HG12	1.95	0.47
1:B:46:LEU:O	1:B:66:ASN:HB3	2.15	0.47
2:D:17:DT:H73	3:E:1:DA:H2	1.79	0.46
1:A:137:ASP:OD1	1:A:139:HIS:N	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:207:ILE:HG13	1:B:233:ILE:HG21	1.97	0.46
1:B:61:LYS:HE3	1:B:84:PHE:HB3	1.97	0.46
1:A:7:HIS:CD2	1:A:238:ARG:CG	2.97	0.46
1:A:18:LEU:HD23	1:A:18:LEU:HA	1.72	0.45
1:A:101:LYS:HE3	1:A:110:THR:OG1	2.15	0.45
1:A:91:ARG:HD3	1:A:93:LYS:HE3	1.99	0.45
1:B:97:TYR:HA	1:B:113:ASN:O	2.17	0.45
2:D:17:DT:H2''	2:D:18:DA:OP2	2.17	0.45
1:B:196:ALA:N	1:B:197:PRO:HD2	2.32	0.45
1:A:83:ASP:OD1	1:A:85:LYS:HG3	2.17	0.44
1:A:152:THR:HG22	1:A:159:GLN:O	2.17	0.44
1:A:126:LYS:HD2	1:A:167:GLU:C	2.38	0.44
1:A:22:ILE:O	1:A:26:ARG:HG3	2.18	0.44
1:A:198:HIS:CD2	1:A:198:HIS:N	2.87	0.43
1:A:221:GLU:HB3	1:A:223:ILE:HD11	2.00	0.43
1:A:101:LYS:HG2	1:A:110:THR:HG23	2.01	0.43
1:A:97:TYR:CE1	1:A:132:LEU:HD11	2.53	0.43
1:A:124:PHE:O	1:A:128:VAL:HG23	2.18	0.43
1:A:127:GLN:O	1:A:127:GLN:HG3	2.19	0.43
1:A:205:ASP:OD2	1:A:230:HIS:HD2	2.02	0.43
1:B:178:ILE:HD12	1:B:183:THR:OG1	2.19	0.43
2:F:2:DC:H2''	2:F:3:DT:H5'	1.99	0.43
1:A:15:ILE:HG21	1:A:18:LEU:HG	2.01	0.42
1:A:182:ARG:HA	1:A:182:ARG:HD3	1.83	0.42
1:A:218:LYS:HE2	1:A:218:LYS:HB3	1.70	0.42
1:A:162:ASN:OD1	1:A:163:LYS:N	2.51	0.42
1:A:194:PHE:CE1	1:A:199:HIS:HA	2.54	0.42
1:B:156:SER:HB3	1:B:201:TYR:HB2	2.00	0.42
1:A:123:ARG:HH11	1:A:123:ARG:HG3	1.83	0.42
1:B:162:ASN:O	1:B:166:GLN:HG3	2.20	0.42
1:B:229:ASP:OD1	1:B:229:ASP:N	2.49	0.41
1:A:23:LYS:HE3	1:A:227:LEU:O	2.21	0.41
1:B:61:LYS:HD2	1:B:63:TYR:OH	2.20	0.41
1:A:129:LEU:HD23	1:A:129:LEU:HA	1.82	0.41
3:E:15:DA:H5'	3:E:15:DA:C8	2.56	0.41
3:G:7:DT:H2''	3:G:8:DA:H5'	2.01	0.41
1:A:146:PHE:HB3	1:A:148:THR:HG22	2.03	0.41
1:A:196:ALA:N	1:A:197:PRO:HD2	2.36	0.41
1:A:20:SER:OG	1:A:22:ILE:HG22	2.20	0.41
1:A:162:ASN:HB3	1:A:165:THR:OG1	2.21	0.41
1:B:84:PHE:HD1	1:B:104:ILE:HG13	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:207:ILE:CD1	1:B:233:ILE:HG22	2.50	0.41
1:A:178:ILE:HD12	1:A:183:THR:OG1	2.21	0.40
1:B:102:GLY:O	1:B:108:GLU:HA	2.21	0.40
1:B:179:ASP:OD2	1:B:179:ASP:C	2.59	0.40
1:B:207:ILE:HD11	1:B:233:ILE:CG2	2.51	0.40
1:A:50:ASP:OD1	1:A:53:ARG:NH1	2.53	0.40
1:A:115:TYR:HD1	1:A:145:ALA:HB3	1.86	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:ARG:NH2	1:A:167:GLU:OE2[8_554]	2.07	0.13

## 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	230/238 (97%)	220 (96%)	10 (4%)	0	100	100
1	B	230/238 (97%)	222 (96%)	8 (4%)	0	100	100
All	All	460/476 (97%)	442 (96%)	18 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	210/215 (98%)	206 (98%)	4 (2%)	57	81
1	B	210/215 (98%)	202 (96%)	8 (4%)	33	64
All	All	420/430 (98%)	408 (97%)	12 (3%)	42	72

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	134	ARG
1	A	156	SER
1	A	174	GLN
1	A	228	SER
1	B	60	ARG
1	B	91	ARG
1	B	134	ARG
1	B	151	SER
1	B	158	ARG
1	B	170	SER
1	B	182	ARG
1	B	187	LYS

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	7	HIS
1	A	106	GLN
1	A	198	HIS
1	B	162	ASN
1	B	166	GLN

### 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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	232/238 (97%)	-0.06	4 (1%) 70 68	18, 41, 69, 81	0
1	B	232/238 (97%)	-0.17	0 100 100	19, 37, 60, 85	0
2	D	18/18 (100%)	0.30	2 (11%) 5 3	39, 53, 114, 128	0
2	F	17/18 (94%)	-0.51	0 100 100	24, 35, 50, 64	0
3	E	17/18 (94%)	0.27	1 (5%) 22 17	29, 51, 109, 117	0
3	G	18/18 (100%)	-0.37	0 100 100	19, 32, 64, 83	0
All	All	534/548 (97%)	-0.11	7 (1%) 77 76	18, 39, 69, 128	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	E	1	DA	4.0
2	D	18	DA	4.0
1	A	7	HIS	3.6
1	A	134	ARG	2.7
1	A	238	ARG	2.5
1	A	216	LYS	2.3
2	D	17	DT	2.3

### 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.