



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 8, 2023 – 10:32 AM EDT

PDB ID : 6ECN  
Title : HIV-1 CA 1/2-hexamer-EE  
Authors : Summers, B.J.; Xiong, Y.  
Deposited on : 2018-08-08  
Resolution : 3.40 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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  
Xtrriage (Phenix) : 1.13  
EDS : 2.35.1  
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) : 2.35.1

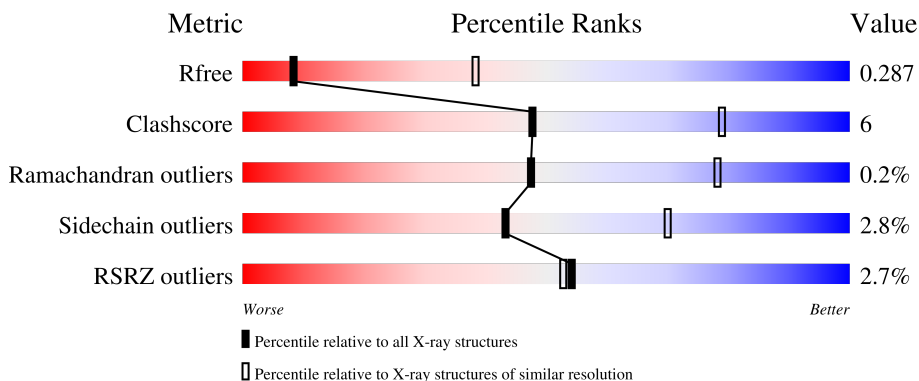
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.40 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	1026 (3.48-3.32)
Clashscore	141614	1055 (3.48-3.32)
Ramachandran outliers	138981	1038 (3.48-3.32)
Sidechain outliers	138945	1038 (3.48-3.32)
RSRZ outliers	127900	2173 (3.50-3.30)

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	231	 71% 14% 13% 2%
1	D	231	 80% 10% 10% 2%
2	B	231	 77% 13% 10% 6%
2	E	231	 66% 19% 14% 3%
3	C	149	 86% 11% 3%

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Mol	Chain	Length	Quality of chain
3	F	149	 80% 14% .. 5%

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8369 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 HIV-1 CA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	201	1539	971	268	288	12	0	0	0
1	D	207	1560	986	268	293	13	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	6	LEU	ILE	conflict	UNP P04591
A	45	CYS	GLU	engineered mutation	UNP P04591
A	54	CYS	THR	engineered mutation	UNP P04591
A	83	LEU	VAL	conflict	UNP P04591
A	120	HIS	ASN	conflict	UNP P04591
A	184	ALA	TRP	engineered mutation	UNP P04591
A	185	ALA	MET	engineered mutation	UNP P04591
A	208	GLY	ALA	conflict	UNP P04591
D	6	LEU	ILE	conflict	UNP P04591
D	45	CYS	GLU	engineered mutation	UNP P04591
D	54	CYS	THR	engineered mutation	UNP P04591
D	83	LEU	VAL	conflict	UNP P04591
D	120	HIS	ASN	conflict	UNP P04591
D	184	ALA	TRP	engineered mutation	UNP P04591
D	185	ALA	MET	engineered mutation	UNP P04591
D	208	GLY	ALA	conflict	UNP P04591

- Molecule 2 is a protein called HIV-1 CA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	207	1535	965	260	298	12	0	0	0
2	E	199	1533	969	263	289	12	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	6	LEU	ILE	conflict	UNP P04591
B	42	CYS	ALA	engineered mutation	UNP P04591
B	54	GLU	THR	engineered mutation	UNP P04591
B	83	LEU	VAL	conflict	UNP P04591
B	120	HIS	ASN	conflict	UNP P04591
B	184	ALA	TRP	engineered mutation	UNP P04591
B	185	ALA	MET	engineered mutation	UNP P04591
B	208	GLY	ALA	conflict	UNP P04591
E	6	LEU	ILE	conflict	UNP P04591
E	42	CYS	ALA	engineered mutation	UNP P04591
E	54	GLU	THR	engineered mutation	UNP P04591
E	83	LEU	VAL	conflict	UNP P04591
E	120	HIS	ASN	conflict	UNP P04591
E	184	ALA	TRP	engineered mutation	UNP P04591
E	185	ALA	MET	engineered mutation	UNP P04591
E	208	GLY	ALA	conflict	UNP P04591

- Molecule 3 is a protein called HIV-1 CA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	145	1113	703	193	208	9	0	0	0
3	F	142	1089	689	188	204	8	0	0	0

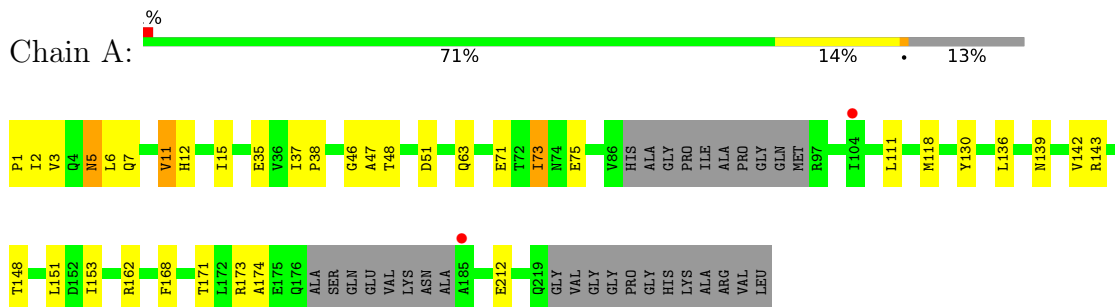
There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	14	CYS	ALA	engineered mutation	UNP A0A248SME6
C	42	GLU	ALA	engineered mutation	UNP A0A248SME6
C	147	LYS	-	expression tag	UNP A0A248SME6
C	148	LEU	-	expression tag	UNP A0A248SME6
C	149	GLN	-	expression tag	UNP A0A248SME6
F	14	CYS	ALA	engineered mutation	UNP A0A248SME6
F	42	GLU	ALA	engineered mutation	UNP A0A248SME6
F	147	LYS	-	expression tag	UNP A0A248SME6
F	148	LEU	-	expression tag	UNP A0A248SME6
F	149	GLN	-	expression tag	UNP A0A248SME6

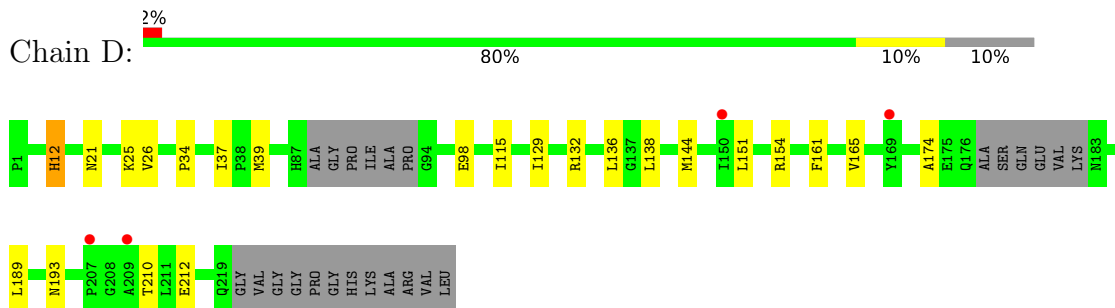
### 3 Residue-property plots [i](#)

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: HIV-1 CA



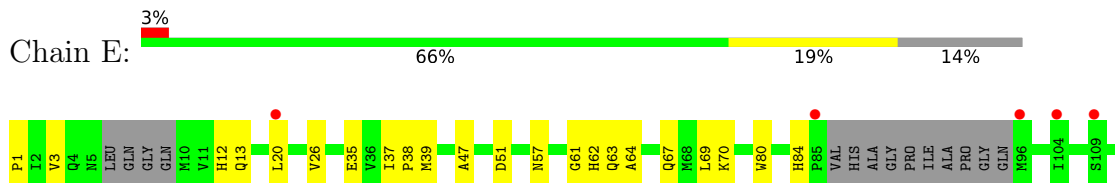
- Molecule 1: HIV-1 CA

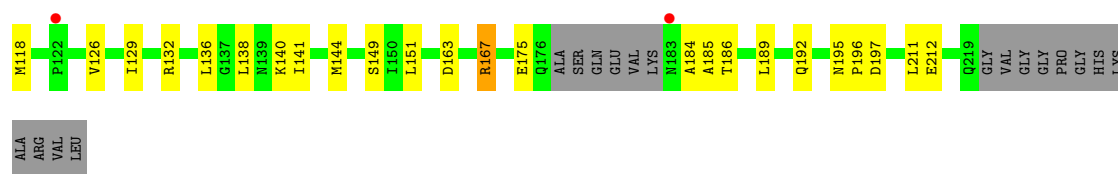


- Molecule 2: HIV-1 CA

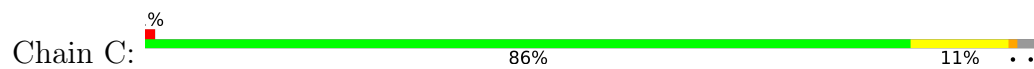


- Molecule 2: HIV-1 CA

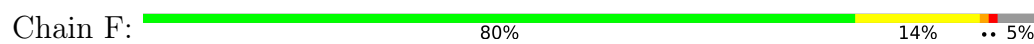




- Molecule 3: HIV-1 CA



- Molecule 3: HIV-1 CA



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.60Å 84.10Å 248.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.01 – 3.40 49.98 – 3.40	Depositor EDS
% Data completeness (in resolution range)	99.7 (50.01-3.40) 99.8 (49.98-3.40)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.70 (at 3.40Å)	Xtrriage
Refinement program	REFMAC 5.8.0103, PHENIX 1.12-2829	Depositor
R, $R_{free}$	0.225 , 0.275 0.237 , 0.287	Depositor DCC
$R_{free}$ test set	955 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	115.1	Xtrriage
Anisotropy	0.226	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.26 , 82.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8369	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	147.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.08% 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 [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	0.58	0/1571	0.77	1/2137 (0.0%)
1	D	0.60	0/1592	0.79	0/2166
2	B	0.59	0/1563	0.77	0/2131
2	E	0.60	0/1564	0.84	3/2125 (0.1%)
3	C	0.66	0/1139	0.79	1/1548 (0.1%)
3	F	0.62	0/1115	0.80	1/1516 (0.1%)
All	All	0.61	0/8544	0.79	6/11623 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	197	ASP	CB-CG-OD2	6.33	124.00	118.30
3	C	136	LEU	CA-CB-CG	-6.28	100.86	115.30
2	E	211	LEU	CB-CG-CD2	-6.04	100.72	111.00
2	E	167	ARG	NE-CZ-NH2	-5.99	117.31	120.30
1	A	173	ARG	NE-CZ-NH1	-5.79	117.41	120.30
3	F	82	ARG	NE-CZ-NH1	5.76	123.18	120.30

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	1539	0	1506	21	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1560	0	1508	14	0
2	B	1535	0	1466	28	0
2	E	1533	0	1501	28	0
3	C	1113	0	1087	10	0
3	F	1089	0	1056	15	0
All	All	8369	0	8124	100	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 6.

All (100) 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)
2:B:69:LEU:HD13	2:B:137:GLY:HA3	1.55	0.87
1:A:3:VAL:HB	1:A:5:ASN:OD1	1.78	0.84
1:D:212:GLU:HG3	3:F:144:MET:HE1	1.64	0.78
1:A:63:GLN:NE2	2:B:173:ARG:HH11	1.83	0.76
2:B:64:ALA:HB2	2:E:63:GLN:CD	2.07	0.75
1:A:63:GLN:HE22	2:B:173:ARG:NH1	1.84	0.75
2:B:62:HIS:O	2:B:63:GLN:NE2	2.23	0.71
1:A:63:GLN:NE2	2:B:173:ARG:NH1	2.41	0.69
2:B:144:MET:HG2	2:E:144:MET:HG2	1.74	0.69
2:E:37:ILE:HD12	2:E:138:LEU:HB3	1.77	0.67
1:D:212:GLU:HG3	3:F:144:MET:CE	2.27	0.64
2:B:69:LEU:HD11	2:B:134:ILE:HA	1.80	0.63
3:F:26:VAL:HG21	3:F:39:MET:HG2	1.80	0.62
1:D:144:MET:HE1	2:E:212:GLU:HG3	1.82	0.61
2:B:65:ALA:HA	2:B:68:MET:CE	2.35	0.57
2:B:188:THR:HG21	2:E:192:GLN:OE1	2.05	0.55
1:D:34:PRO:HG3	1:D:174:ALA:HA	1.88	0.54
2:B:65:ALA:HA	2:B:68:MET:HE3	1.90	0.54
1:A:6:LEU:CB	1:A:11:VAL:HG22	2.37	0.54
2:B:69:LEU:HD21	2:B:134:ILE:HG23	1.90	0.54
2:B:144:MET:HG2	2:E:144:MET:CG	2.38	0.54
3:C:37:ILE:HD12	3:C:138:LEU:HB3	1.89	0.54
1:D:37:ILE:HD12	1:D:138:LEU:HB3	1.88	0.53
1:A:12:HIS:ND1	1:A:111:LEU:HD11	2.22	0.53
2:E:1:PRO:H2	2:E:13:GLN:HB2	1.74	0.53
2:B:37:ILE:HD12	2:B:138:LEU:HB3	1.91	0.53
1:A:212:GLU:OE1	3:C:140:LYS:HE2	2.08	0.53
2:B:144:MET:CG	2:E:144:MET:HG2	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:81:ASP:OD1	2:B:100:ARG:HG2	2.09	0.52
2:E:163:ASP:O	2:E:167:ARG:HG3	2.10	0.52
2:E:184:ALA:C	2:E:186:THR:H	2.13	0.52
3:F:79:GLU:OE1	3:F:82:ARG:NH2	2.43	0.52
2:E:26:VAL:HG21	2:E:39:MET:HG2	1.90	0.52
2:E:184:ALA:O	2:E:185:ALA:HB3	2.10	0.52
1:A:1:PRO:HG2	1:A:15:ILE:HA	1.91	0.52
1:D:151:LEU:HD23	1:D:189:LEU:HD11	1.91	0.51
2:E:140:LYS:O	2:E:144:MET:HB2	2.10	0.51
3:C:37:ILE:HB	3:C:38:PRO:HD3	1.93	0.50
3:F:4:GLN:HG2	3:F:10:MET:SD	2.51	0.50
3:F:37:ILE:CG2	3:F:135:ILE:HG23	2.42	0.50
2:E:80:TRP:O	2:E:84:HIS:N	2.44	0.50
1:A:37:ILE:HD11	1:A:142:VAL:HG21	1.94	0.49
2:B:163:ASP:O	2:B:167:ARG:HG3	2.12	0.49
1:D:129:ILE:HG12	1:D:132:ARG:HH12	1.77	0.48
1:A:171:THR:O	1:A:174:ALA:HB3	2.13	0.48
2:B:97:ARG:NH1	2:B:103:ASP:OD2	2.43	0.48
1:A:153:ILE:HG21	1:A:168:PHE:HA	1.95	0.48
1:D:26:VAL:HG21	1:D:39:MET:HG2	1.95	0.48
3:C:37:ILE:CD1	3:C:138:LEU:HB3	2.44	0.48
1:D:37:ILE:CD1	1:D:138:LEU:HB3	2.43	0.48
2:E:47:ALA:HB1	2:E:51:ASP:HB2	1.95	0.48
2:E:195:ASN:HB2	2:E:196:PRO:HD2	1.94	0.48
2:E:149:SER:HA	2:E:175:GLU:OE2	2.14	0.48
1:D:21:ASN:O	1:D:25:LYS:HG2	2.13	0.48
2:B:69:LEU:CD1	2:B:137:GLY:HA3	2.37	0.48
2:B:189:LEU:HD22	2:E:151:LEU:HD21	1.96	0.47
1:A:47:ALA:HB1	1:A:51:ASP:HB2	1.95	0.47
3:C:79:GLU:OE2	3:C:83:LEU:HD11	2.14	0.47
3:F:26:VAL:O	3:F:30:LYS:HB3	2.14	0.47
3:F:72:THR:HG21	3:F:137:GLY:HA2	1.96	0.47
1:D:161:PHE:O	1:D:165:VAL:HG23	2.15	0.47
3:F:82:ARG:HH11	3:F:82:ARG:HB3	1.80	0.47
2:B:64:ALA:HB3	2:E:64:ALA:HB2	1.98	0.46
1:A:35:GLU:O	1:A:38:PRO:HD2	2.15	0.46
1:A:73:ILE:HG12	1:A:130:TYR:HE1	1.79	0.46
3:C:97:ARG:NH2	3:C:103:ASP:OD2	2.48	0.46
2:E:61:GLY:O	2:E:62:HIS:C	2.54	0.46
2:B:35:GLU:O	2:B:38:PRO:HD2	2.16	0.46
2:B:187:GLU:O	2:B:191:VAL:HG23	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:12:HIS:HB3	1:D:115:ILE:HD11	1.97	0.45
1:A:5:ASN:OD1	1:A:5:ASN:N	2.46	0.45
1:A:37:ILE:HB	1:A:38:PRO:HD3	1.98	0.45
2:B:64:ALA:O	2:B:68:MET:N	2.49	0.45
3:F:37:ILE:HG21	3:F:135:ILE:HG23	1.99	0.45
1:A:2:ILE:HG22	1:A:46:GLY:O	2.17	0.44
2:B:64:ALA:HB2	2:E:63:GLN:OE1	2.16	0.44
2:B:64:ALA:O	2:B:67:GLN:N	2.51	0.44
2:E:20:LEU:HD23	2:E:20:LEU:HA	1.93	0.44
1:A:48:THR:HG22	1:A:118:MET:SD	2.58	0.44
3:F:30:LYS:HD2	3:F:33:SER:HB3	1.99	0.44
3:C:136:LEU:HD23	3:C:136:LEU:HA	1.81	0.44
2:E:129:ILE:HG12	2:E:132:ARG:HH12	1.82	0.43
3:F:23:TRP:O	3:F:27:VAL:HG23	2.18	0.43
3:C:23:TRP:NE1	3:C:55:MET:O	2.50	0.43
2:E:69:LEU:HB2	2:E:141:ILE:HD11	2.00	0.42
2:B:153:ILE:HG21	2:B:168:PHE:HA	2.01	0.42
3:C:54:THR:O	3:C:58:THR:HG23	2.19	0.42
2:B:65:ALA:HA	2:B:68:MET:HE2	2.02	0.42
2:E:189:LEU:HA	2:E:189:LEU:HD12	1.72	0.42
1:A:71:GLU:O	1:A:75:GLU:HG3	2.21	0.41
2:E:35:GLU:O	2:E:38:PRO:HD2	2.21	0.41
1:D:154:ARG:HA	1:D:193:ASN:HB3	2.02	0.41
2:E:118:MET:HG3	2:E:126:VAL:HG22	2.03	0.41
1:A:35:GLU:C	1:A:38:PRO:HD2	2.41	0.40
3:C:21:ASN:O	3:C:25:LYS:HG2	2.21	0.40
1:D:210:THR:HB	3:F:71:GLU:OE1	2.21	0.40
3:F:118:MET:HG3	3:F:126:VAL:HG22	2.02	0.40
2:E:3:VAL:HB	2:E:13:GLN:HG3	2.03	0.40
3:F:33:SER:OG	3:F:34:PRO:HD2	2.21	0.40
1:A:139:ASN:O	1:A:143:ARG:HG3	2.21	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	195/231 (84%)	188 (96%)	6 (3%)	1 (0%)	29	61
1	D	201/231 (87%)	196 (98%)	5 (2%)	0	100	100
2	B	201/231 (87%)	196 (98%)	4 (2%)	1 (0%)	29	61
2	E	191/231 (83%)	182 (95%)	9 (5%)	0	100	100
3	C	141/149 (95%)	140 (99%)	1 (1%)	0	100	100
3	F	138/149 (93%)	135 (98%)	3 (2%)	0	100	100
All	All	1067/1222 (87%)	1037 (97%)	28 (3%)	2 (0%)	47	78

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
2	B	8	GLY

### 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	163/192 (85%)	156 (96%)	7 (4%)	29	59
1	D	161/192 (84%)	158 (98%)	3 (2%)	57	78
2	B	157/193 (81%)	155 (99%)	2 (1%)	69	84
2	E	163/193 (84%)	158 (97%)	5 (3%)	40	68
3	C	119/128 (93%)	117 (98%)	2 (2%)	60	80
3	F	116/128 (91%)	110 (95%)	6 (5%)	23	53
All	All	879/1026 (86%)	854 (97%)	25 (3%)	43	70

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

Mol	Chain	Res	Type
1	A	5	ASN
1	A	11	VAL
1	A	73	ILE
1	A	136	LEU
1	A	148	THR
1	A	151	LEU
1	A	162	ARG
2	B	3	VAL
2	B	136	LEU
3	C	18	ARG
3	C	33	SER
1	D	12	HIS
1	D	98	GLU
1	D	136	LEU
2	E	12	HIS
2	E	57	ASN
2	E	67	GLN
2	E	70	LYS
2	E	136	LEU
3	F	4	GLN
3	F	18	ARG
3	F	82	ARG
3	F	87	HIS
3	F	125	PRO
3	F	136	LEU

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

Mol	Chain	Res	Type
1	A	63	GLN
2	B	63	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 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

### 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<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	201/231 (87%)	-0.04	2 (0%) 82 81	94, 140, 223, 277	0
1	D	207/231 (89%)	-0.14	4 (1%) 66 65	77, 130, 183, 214	0
2	B	207/231 (89%)	0.39	15 (7%) 15 17	90, 181, 251, 297	0
2	E	199/231 (86%)	0.16	7 (3%) 44 43	86, 149, 239, 273	0
3	C	145/149 (97%)	0.00	2 (1%) 75 74	82, 117, 176, 226	0
3	F	142/149 (95%)	-0.15	0 100 100	93, 133, 197, 230	0
All	All	1101/1222 (90%)	0.05	30 (2%) 54 53	77, 139, 228, 297	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	104	ILE	5.4
2	E	85	PRO	4.8
2	E	183	ASN	4.6
2	B	47	ALA	4.3
2	B	101	GLY	4.1
1	A	185	ALA	3.9
2	B	46	GLY	3.8
2	B	118	MET	3.8
2	B	8	GLY	3.6
3	C	4	GLN	3.6
1	D	207	PRO	3.5
2	B	128	GLU	3.4
2	B	117	TRP	3.3
1	D	209	ALA	3.2
2	E	104	ILE	3.0
1	D	150	ILE	2.9
3	C	115	ILE	2.6
2	B	7	GLN	2.6
2	B	121	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
2	E	109	SER	2.5
2	B	10	MET	2.4
2	E	96	MET	2.4
2	E	20	LEU	2.4
2	B	122	PRO	2.3
2	E	122	PRO	2.3
2	B	183	ASN	2.3
1	A	104	ILE	2.3
2	B	119	THR	2.2
1	D	169	TYR	2.1
2	B	115	ILE	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 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.