



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:06 PM UTC

PDB ID : 6EC2 / pdb_00006ec2
Title : Structure of HIV-1 CA 1/3-hexamer
Authors : Summers, B.J.; Xiong, Y.
Deposited on : 2018-08-07
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

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>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

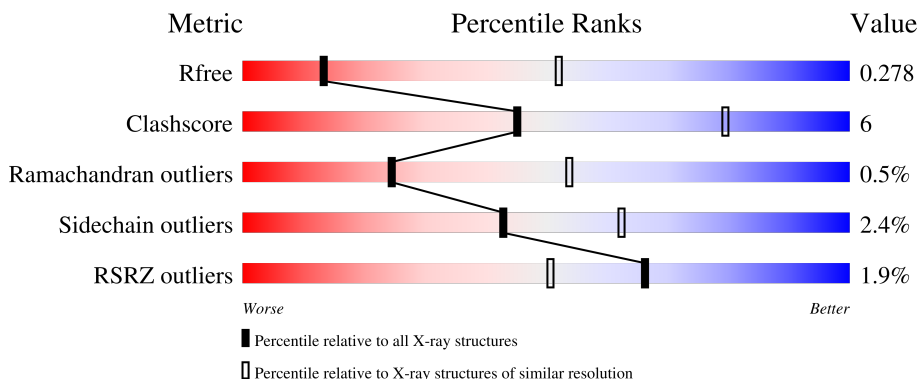
1 Overall quality at a glance i

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}	180053	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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 $\leq 5\%$. 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	 74% 15% • 8%
1	F	231	 75% 16% • 8%
2	C	231	 78% 8% • 13%
2	G	231	 81% 6% 13%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6405 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 Capsid protein p24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	212	1633	1029	280	311	13	0	9	0
1	F	212	1639	1031	286	309	13	0	10	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	GLU	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
F	6	LEU	ILE	conflict	UNP P04591
F	45	CYS	GLU	engineered mutation	UNP P04591
F	54	GLU	THR	engineered mutation	UNP P04591
F	83	LEU	VAL	conflict	UNP P04591
F	120	HIS	ASN	conflict	UNP P04591
F	184	ALA	TRP	engineered mutation	UNP P04591
F	185	ALA	MET	engineered mutation	UNP P04591
F	208	GLY	ALA	conflict	UNP P04591

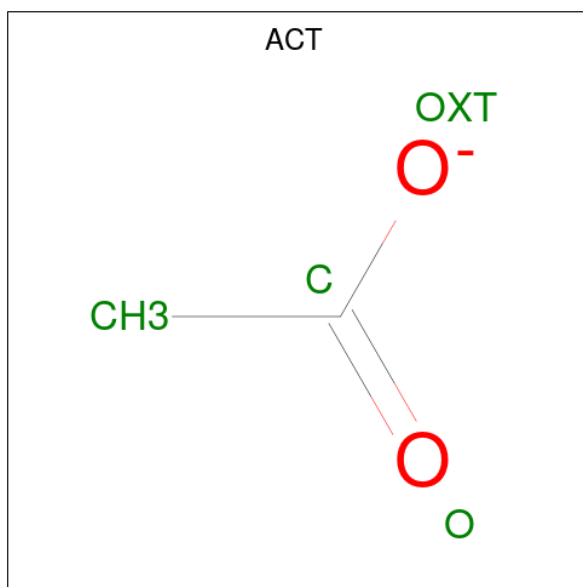
- Molecule 2 is a protein called Capsid protein p24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	201	1568	993	267	295	13	0	9	0
2	G	201	1539	977	262	288	12	0	9	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	6	LEU	ILE	conflict	UNP P04591
C	14	CYS	ALA	engineered mutation	UNP P04591
C	42	GLU	ALA	engineered mutation	UNP P04591
C	83	LEU	VAL	conflict	UNP P04591
C	120	HIS	ASN	conflict	UNP P04591
C	184	ALA	TRP	engineered mutation	UNP P04591
C	185	ALA	MET	engineered mutation	UNP P04591
C	208	GLY	ALA	conflict	UNP P04591
G	6	LEU	ILE	conflict	UNP P04591
G	14	CYS	ALA	engineered mutation	UNP P04591
G	42	GLU	ALA	engineered mutation	UNP P04591
G	83	LEU	VAL	conflict	UNP P04591
G	120	HIS	ASN	conflict	UNP P04591
G	184	ALA	TRP	engineered mutation	UNP P04591
G	185	ALA	MET	engineered mutation	UNP P04591
G	208	GLY	ALA	conflict	UNP P04591

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf		
3	A	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	F	1	Total	C	O	0	0
			4	2	2		

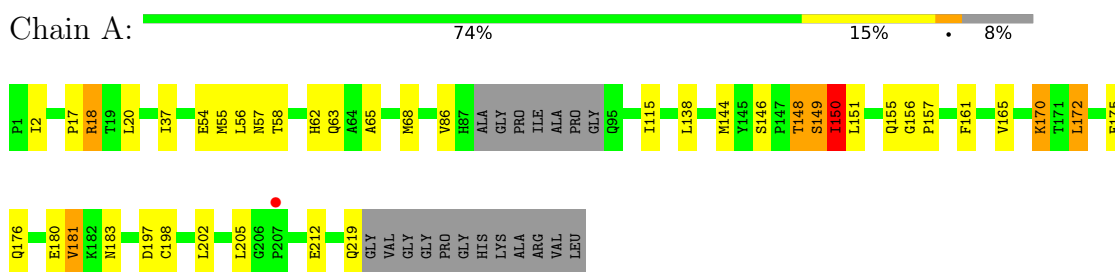
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	4	Total O 4 4	0	0
4	C	4	Total O 4 4	0	0
4	F	6	Total O 6 6	0	0

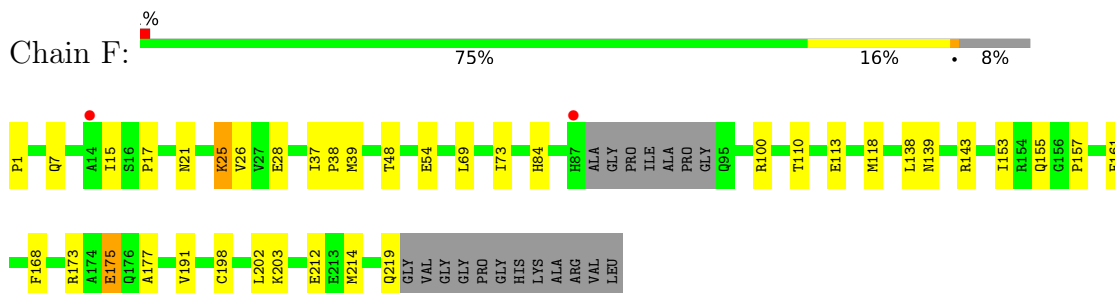
3 Residue-property plots [i](#)

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.

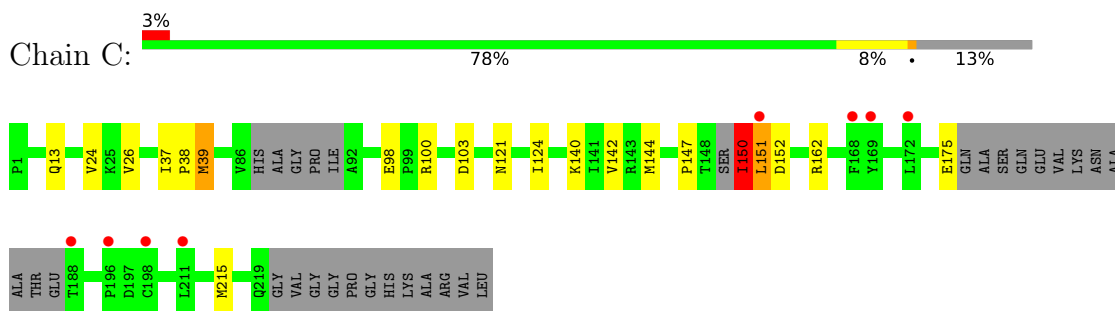
- Molecule 1: Capsid protein p24



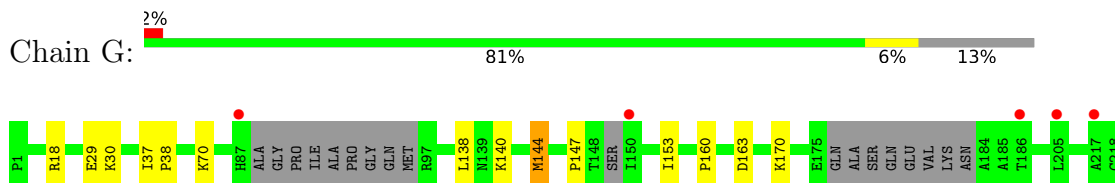
- Molecule 1: Capsid protein p24



- Molecule 2: Capsid protein p24



- Molecule 2: Capsid protein p24



0219
GLY
VAL
GLY
GLY
PRO
GLY
HIS
LYS
ALA
ARG
VAL
LEU

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.80Å 97.47Å 150.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.01 – 3.40 50.01 – 3.40	Depositor EDS
% Data completeness (in resolution range)	99.3 (50.01-3.40) 99.3 (50.01-3.40)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 3.41Å)	Xtrriage
Refinement program	REFMAC 5.8.0103, PHENIX 1.12_2829	Depositor
R, R_{free}	0.211 , 0.275 0.213 , 0.278	Depositor DCC
R_{free} test set	806 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	66.0	Xtrriage
Anisotropy	0.810	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 64.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6405	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.84% 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

Bond lengths and bond angles in the following residue types are not validated in this section: ACT

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	1.10	6/1669 (0.4%)	1.28	14/2271 (0.6%)
1	F	1.02	1/1675 (0.1%)	1.10	2/2277 (0.1%)
2	C	0.97	2/1603 (0.1%)	1.07	2/2177 (0.1%)
2	G	0.86	0/1573	1.03	0/2139
All	All	0.99	9/6520 (0.1%)	1.12	18/8864 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	150[A]	ILE	C-O	8.16	1.33	1.24
1	A	55	MET	C-O	-6.36	1.16	1.24
1	A	149	SER	CA-C	-6.08	1.45	1.52
2	C	124	ILE	CA-C	5.74	1.58	1.52
2	C	13	GLN	CA-C	-5.65	1.45	1.52
1	F	28	GLU	C-O	5.51	1.30	1.24
1	A	54	GLU	C-O	-5.21	1.18	1.24
1	A	170	LYS	C-O	5.15	1.30	1.24
1	A	197	ASP	C-O	5.14	1.30	1.24

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	150[A]	ILE	CA-C-O	10.06	133.36	120.78
1	A	150[A]	ILE	N-CA-CB	-7.84	98.30	111.23
1	A	155	GLN	N-CA-C	7.20	119.12	111.28
1	A	18	ARG	CB-CA-C	-6.81	100.18	110.88
2	C	39	MET	CG-SD-CE	6.31	114.78	100.90
1	A	197	ASP	CB-CA-C	-6.18	101.09	110.92
1	A	63	GLN	CB-CA-C	-5.94	99.64	110.63
1	A	62	HIS	CB-CA-C	-5.74	103.42	111.80
1	A	150[A]	ILE	N-CA-C	5.71	121.21	109.34
1	A	170	LYS	N-CA-CB	5.46	117.98	110.07
1	A	172	LEU	N-CA-C	-5.28	106.10	112.54
2	C	150[A]	ILE	CA-C-O	5.26	129.74	120.80
1	F	139	ASN	CB-CA-C	-5.19	102.18	110.79
1	A	56	LEU	CA-C-O	-5.14	115.11	120.55
1	A	57	ASN	N-CA-C	5.07	118.56	112.38
1	A	86	VAL	N-CA-C	5.04	115.83	108.58
1	A	170	LYS	CB-CG-CD	5.02	122.85	111.30
1	F	25	LYS	CB-CA-C	-5.02	102.99	110.88

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	156	GLY	Peptide

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	1633	0	1592	24	0
1	F	1639	0	1602	26	0
2	C	1568	0	1532	19	0
2	G	1539	0	1498	14	0
3	A	4	0	3	0	0
3	C	4	0	3	0	0
3	F	4	0	3	0	0
4	A	4	0	0	0	0
4	C	4	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	F	6	0	0	1	0
All	All	6405	0	6233	74	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 (74) 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:F:15:ILE:CG2	1:F:54:GLU:OE1	1.87	1.22
1:A:146:SER:OG	1:A:175:GLU:HG2	1.37	1.19
1:A:148:THR:O	1:A:175:GLU:HG3	1.57	1.04
1:F:15:ILE:HG22	1:F:54:GLU:OE1	1.55	1.04
1:F:15:ILE:HG21	1:F:54:GLU:OE1	1.57	1.03
2:C:150[A]:ILE:HG23	2:C:151:LEU:N	1.78	0.97
2:C:147:PRO:HB2	2:C:175:GLU:OE2	1.64	0.95
2:C:150[A]:ILE:CG2	2:C:151:LEU:H	1.79	0.95
1:A:68:MET:SD	1:A:144:MET:HE1	2.12	0.90
1:A:148:THR:O	1:A:175:GLU:CG	2.21	0.88
1:F:84:HIS:O	1:F:100:ARG:NH1	2.09	0.85
2:C:150[A]:ILE:HG23	2:C:151:LEU:H	1.37	0.81
1:F:157[A]:PRO:HD2	4:F:402:HOH:O	1.83	0.78
1:A:146:SER:HG	1:A:175:GLU:HG2	1.52	0.73
1:A:146:SER:OG	1:A:175:GLU:CG	2.29	0.72
1:A:172:LEU:O	1:A:176:GLN:HB3	1.91	0.70
1:A:37:ILE:HD12	1:A:138:LEU:HB3	1.79	0.65
1:A:180:GLU:HG2	2:G:70:LYS:HE2	1.78	0.64
1:A:212:GLU:HG3	2:G:144:MET:HE1	1.80	0.64
2:G:30:LYS:HE2	2:G:163:ASP:OD2	2.02	0.59
2:G:37:ILE:HD12	2:G:138:LEU:HB3	1.84	0.59
2:C:144:MET:CE	1:F:212:GLU:HG3	2.33	0.58
1:A:144:MET:O	1:F:143:ARG:NH1	2.37	0.57
2:C:26[A]:VAL:HG21	2:C:39:MET:HE2	1.85	0.57
2:G:147:PRO:HG2	2:G:153:ILE:CD1	2.34	0.57
1:A:65:ALA:HA	1:A:68:MET:HE3	1.87	0.56
1:A:2:ILE:HD12	1:A:115:ILE:HG23	1.88	0.55
1:A:148:THR:O	1:A:175:GLU:HG2	2.07	0.54
1:A:149:SER:O	1:A:151:LEU:N	2.40	0.54
2:C:26[B]:VAL:HG11	2:C:39:MET:HE2	1.89	0.53
2:C:147:PRO:CB	2:C:175:GLU:OE2	2.48	0.53
1:A:18:ARG:CZ	2:G:18:ARG:HD2	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:144:MET:HE2	1:F:212:GLU:HG3	1.91	0.52
2:G:29:GLU:HB3	2:G:160:PRO:HD3	1.92	0.52
1:F:155:GLN:HE22	1:F:161:PHE:HD1	1.58	0.52
2:G:30:LYS:HE2	2:G:163:ASP:CG	2.35	0.52
1:F:37:ILE:HD12	1:F:138:LEU:HB3	1.92	0.51
1:F:26[A]:VAL:HG21	1:F:39:MET:HG2	1.93	0.51
1:F:175:GLU:O	1:F:177:ALA:N	2.43	0.50
2:G:37:ILE:HB	2:G:38[A]:PRO:HD3	1.92	0.50
1:A:150[A]:ILE:HG23	1:A:151:LEU:HD23	1.91	0.50
2:G:147:PRO:HG2	2:G:153:ILE:HD11	1.94	0.49
2:C:150[A]:ILE:HG22	2:C:151:LEU:H	1.75	0.48
1:A:20:LEU:HD22	1:A:58:THR:HG21	1.97	0.47
1:A:172:LEU:O	1:A:176:GLN:CB	2.61	0.46
4:C:401:HOH:O	1:F:173:ARG:HD3	2.16	0.45
2:G:170:LYS:HD2	2:G:170:LYS:HA	1.73	0.45
1:F:37:ILE:HB	1:F:38[A]:PRO:HD3	1.97	0.45
1:F:191:VAL:HG22	1:F:214[A]:MET:HE1	1.98	0.45
2:G:147:PRO:HG2	2:G:153:ILE:HD13	1.98	0.44
2:C:140:LYS:HE3	1:F:212:GLU:OE1	2.17	0.44
2:G:37:ILE:CD1	2:G:138:LEU:HB3	2.47	0.44
1:A:172:LEU:HG	1:A:181:VAL:HG11	1.99	0.44
1:A:149:SER:C	1:A:151:LEU:N	2.76	0.44
2:C:144:MET:HE1	1:F:212:GLU:HG3	1.98	0.44
1:F:110:THR:OG1	1:F:113:GLU:HG3	2.18	0.43
1:F:191:VAL:CG2	1:F:214[A]:MET:HE1	2.48	0.43
1:F:153:ILE:HG21	1:F:168:PHE:HA	2.02	0.42
2:C:162:ARG:HA	2:C:215:MET:HE3	2.02	0.42
2:C:37:ILE:HB	2:C:38:PRO:HD3	2.01	0.42
1:F:69:LEU:O	1:F:73:ILE:HG13	2.19	0.42
1:A:202:LEU:O	1:A:205:LEU:HB2	2.20	0.42
1:F:21:ASN:O	1:F:25:LYS:HG2	2.20	0.42
1:A:161:PHE:O	1:A:165:VAL:HG23	2.20	0.41
2:C:37:ILE:HD11	2:C:142:VAL:HG21	2.02	0.41
2:C:100:ARG:HG2	2:C:103:ASP:OD2	2.21	0.41
1:F:1[A]:PRO:HD3	1:F:15:ILE:HA	2.02	0.41
1:F:7:GLN:OE1	1:F:7:GLN:N	2.53	0.41
1:A:212:GLU:OE2	2:G:140:LYS:HE3	2.21	0.41
1:F:48:THR:HG22	1:F:118:MET:SD	2.61	0.41
1:F:202:LEU:HD22	1:F:214[A]:MET:HB3	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	209/231 (90%)	203 (97%)	3 (1%)	3 (1%)	9	31
1	F	209/231 (90%)	199 (95%)	10 (5%)	0	100	100
2	C	195/231 (84%)	193 (99%)	1 (0%)	1 (0%)	24	54
2	G	194/231 (84%)	192 (99%)	2 (1%)	0	100	100
All	All	807/924 (87%)	787 (98%)	16 (2%)	4 (0%)	24	54

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	181	VAL
1	A	157[A]	PRO
2	C	151	LEU
1	A	150[A]	ILE

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	174/192 (91%)	168 (97%)	6 (3%)	32	57
1	F	175/192 (91%)	170 (97%)	5 (3%)	37	60
2	C	169/194 (87%)	165 (98%)	4 (2%)	43	62
2	G	162/194 (84%)	161 (99%)	1 (1%)	78	80
All	All	680/772 (88%)	664 (98%)	16 (2%)	43	62

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	PRO
1	A	148	THR
1	A	170	LYS
1	A	183	ASN
1	A	198	CYS
1	A	219	GLN
2	C	24	VAL
2	C	98	GLU
2	C	121	ASN
2	C	150[A]	ILE
1	F	17	PRO
1	F	175	GLU
1	F	198	CYS
1	F	203	LYS
1	F	219	GLN
2	G	144	MET

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	155	GLN
1	A	176	GLN
1	A	183	ASN
2	C	50	GLN
2	C	84	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	ACT	F	301	-	3,3,3	1.10	0	3,3,3	0.83	0
3	ACT	C	301	-	3,3,3	0.86	0	3,3,3	0.90	0
3	ACT	A	301	-	3,3,3	0.86	0	3,3,3	1.41	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	212/231 (91%)	-0.29	1 (0%) 87 78	37, 58, 91, 112	2 (0%)
1	F	212/231 (91%)	-0.26	2 (0%) 81 68	34, 59, 100, 133	2 (0%)
2	C	201/231 (87%)	0.08	8 (3%) 42 30	21, 57, 175, 215	4 (1%)
2	G	201/231 (87%)	0.16	5 (2%) 58 43	46, 83, 166, 213	3 (1%)
All	All	826/924 (89%)	-0.08	16 (1%) 66 51	21, 62, 160, 215	11 (1%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	G	87	HIS	3.6
2	G	186	THR	3.2
2	C	198	CYS	3.0
2	G	205	LEU	2.8
2	G	150[A]	ILE	2.8
2	C	168	PHE	2.7
2	C	211	LEU	2.6
2	G	217	ALA	2.4
1	F	87	HIS	2.4
2	C	151	LEU	2.4
2	C	196[A]	PRO	2.4
2	C	188	THR	2.3
2	C	169	TYR	2.3
2	C	172	LEU	2.2
1	F	14	ALA	2.1
1	A	207	PRO	2.1

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	ACT	F	301	4/4	0.75	0.13	59,63,63,64	0
3	ACT	C	301	4/4	0.77	0.14	89,95,98,100	0
3	ACT	A	301	4/4	0.88	0.20	69,76,76,76	0

6.5 Other polymers [i](#)

There are no such residues in this entry.