

Full wwPDB X-ray Structure Validation Report (i)

Aug 20, 2023 - 09:06 AM EDT

:	2ONT
:	A swapped dimer of the HIV-1 capsid C-terminal domain
:	Ivanov, D.; Tsodikov, O.V.; Kasanov, J.; Ellenberger, T.; Wagner, G.; Collins,
	Т.
	2007-01-24
:	2.40 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 (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 (1)) were used in the production of this report:

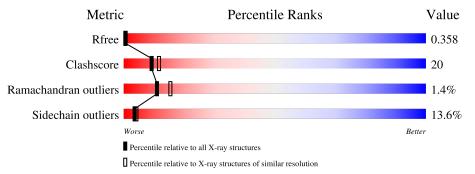
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)		
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.40 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)	$\begin{array}{c} {\rm Similar\ resolution} \\ (\# {\rm Entries,\ resolution\ range}({\rm \AA})) \end{array}$
R _{free}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)

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%

Mol	Chain	Length	Quality of chain						
1	А	76	62%	28%	7% •				



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 580 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		
1	А	73	Total 561	C 352	N 96	0 108	${ m S}{ m 5}$	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	145	GLY	-	cloning artifact	UNP P12497
А	?	-	ALA	deletion	UNP P12497

• Molecule 2 is water.

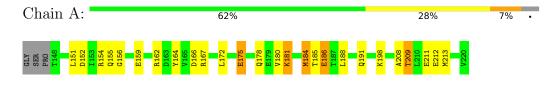
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	19	Total O 19 19	0	0



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. 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. 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





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants	40.13Å 40.13Å 105.32Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 - 2.40	Depositor
Resolution (A)	31.92 - 2.30	EDS
% Data completeness	99.2 (30.00-2.40)	Depositor
(in resolution range)	98.2 (31.92-2.30)	EDS
R _{merge}	0.05	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$4.24 (at 2.31 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
D D	0.243 , 0.294	Depositor
R, R_{free}	0.300 , 0.358	DCC
R_{free} test set	198 reflections (4.78%)	wwPDB-VP
Wilson B-factor $(Å^2)$	52.6	Xtriage
Anisotropy	0.356	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.34, 26.2	EDS
L-test for twinning ²	$ \langle L \rangle = 0.47, \langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	580	wwPDB-VP
Average B, all atoms $(Å^2)$	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 10.92% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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	Bon	d lengths	Bond angles		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.89	1/570~(0.2%)	0.83	2/773~(0.3%)	

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
1	А	186	GLU	CG-CD	5.33	1.59	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
1	А	184	MET	CB-CG-SD	6.20	130.99	112.40
1	А	166	ASP	CB-CG-OD1	5.01	122.81	118.30

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	А	561	0	544	22	1
2	А	19	0	0	8	0
All	All	580	0	544	22	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 20.

All (22) 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:151:LEU:HB3	2:A:15:HOH:O	1.63	0.97
1:A:181:LYS:O	1:A:185:THR:HG23	1.72	0.88
1:A:209:THR:HG22	1:A:212:GLU:H	1.37	0.88
1:A:175:GLU:HG2	1:A:180:VAL:CG1	2.19	0.73
1:A:151:LEU:HG	2:A:16:HOH:O	1.93	0.69
1:A:211:GLU:HB2	2:A:3:HOH:O	1.92	0.69
1:A:209:THR:O	1:A:213:MET:HG3	2.01	0.60
1:A:209:THR:HG22	1:A:212:GLU:N	2.16	0.55
1:A:162:ARG:HG3	2:A:8:HOH:O	2.06	0.54
1:A:159:GLU:O	2:A:14:HOH:O	2.18	0.53
1:A:155:GLN:HG2	1:A:156:GLY:O	2.10	0.52
1:A:208:ALA:HB3	2:A:4:HOH:O	2.08	0.52
1:A:155:GLN:HB2	1:A:164:TYR:CG	2.45	0.51
1:A:154:ARG:O	1:A:167:ARG:NH2	2.44	0.49
1:A:175:GLU:HG2	1:A:180:VAL:HG12	1.94	0.49
1:A:209:THR:CG2	1:A:211:GLU:HB3	2.44	0.48
1:A:155:GLN:HB2	1:A:164:TYR:CD1	2.48	0.48
1:A:211:GLU:HG2	2:A:19:HOH:O	2.15	0.46
1:A:209:THR:HG21	1:A:211:GLU:HB3	1.99	0.45
1:A:154:ARG:NE	2:A:17:HOH:O	2.51	0.43
1:A:181:LYS:O	1:A:184:MET:HB3	2.17	0.43
1:A:175:GLU:CG	1:A:180:VAL:CG1	2.94	0.41

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:152:ASP:OD1	1:A:162:ARG:NH2[3_554]	2.14	0.06

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	А	71/76~(93%)	68~(96%)	2(3%)	1 (1%)	11 15

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	198	LYS

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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	А	59/66~(89%)	51 (86%)	8 (14%)	3 4

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	172	LEU
1	А	175	GLU
1	А	178	GLN
1	А	181	LYS
1	А	186	GLU
1	А	188	LEU
1	А	191	GLN
1	А	209	THR

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

Mol	Chain	Res	Type
1	А	155	GLN
1	А	178	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 (i)

6.1 Protein, DNA and RNA chains (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

