



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 09:07 PM GMT

PDB ID : 1NJU
Title : Complex structure of HCMV Protease and a peptidomimetic inhibitor
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Deposited on : 2003-01-02
Resolution : 2.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

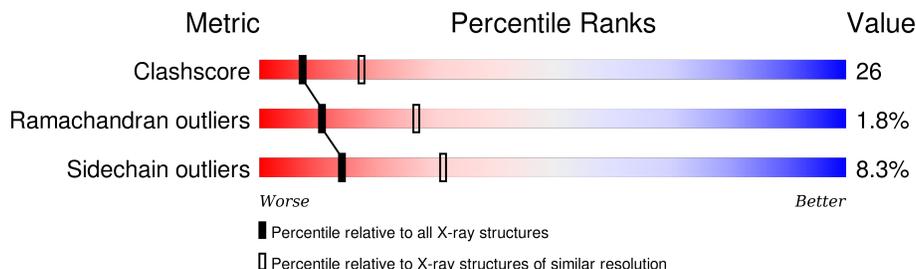
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	256	
1	B	256	
1	C	256	
1	D	256	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7392 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.

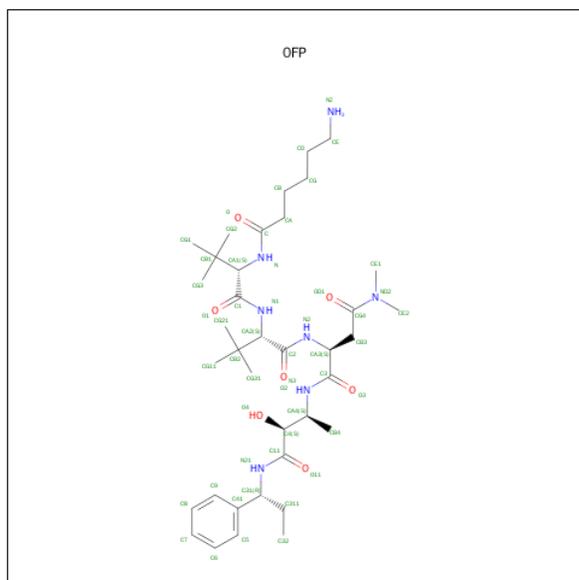
- Molecule 1 is a protein called Assemblin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	230	1804	1128	327	344	5	0	0	0
1	B	230	1801	1125	327	344	5	0	0	0
1	C	227	1779	1113	323	338	5	0	0	0
1	D	230	1804	1128	327	344	5	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	143	GLN	ALA	ENGINEERED	UNP P16753
B	443	GLN	ALA	ENGINEERED	UNP P16753
C	1143	GLN	ALA	ENGINEERED	UNP P16753
D	1443	GLN	ALA	ENGINEERED	UNP P16753

- Molecule 2 is N-(6-AMINOHEXANOYL)-3-METHYL-L-VALYL-3-METHYL-L-VALYL-N 1 -[(2S,3S)-3-HYDROXY-4-OXO-4-[(1R)-1-PHENYLPROPYL]AMINO}BUTAN-2-YL]-N 4 ,N 4 -DIMETHYL-L-ASPARTAMIDE (three-letter code: 0FP) (formula: C₃₇H₆₃N₇O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	N			O
2	A	1	51	37	7	7	0	0
2	B	1	51	37	7	7	0	0
2	C	1	51	37	7	7	0	0
2	D	1	51	37	7	7	0	0



- Molecule 1: Assemblin

Chain D: 51% 32% 7% 10%



4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	105.46Å 213.53Å 52.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.88 – 2.70	Depositor
% Data completeness (in resolution range)	81.7 (25.88-2.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.223 , 0.273	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	7392	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 0FP

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	0.40	0/1839	0.63	0/2492
1	B	0.41	0/1836	0.64	0/2488
1	C	0.39	0/1814	0.62	0/2458
1	D	0.39	0/1839	0.63	0/2492
All	All	0.40	0/7328	0.63	0/9930

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	D	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	1530	TYR	Sidechain

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	1804	0	1768	88	0
1	B	1801	0	1762	96	0
1	C	1779	0	1745	104	0
1	D	1804	0	1768	110	0
2	A	51	0	62	8	0
2	B	51	0	62	6	0
2	C	51	0	62	14	0
2	D	51	0	62	2	0
All	All	7392	0	7291	379	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 26.

The worst 5 of 379 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:560:0FP:H	2:B:560:0FP:HD1	1.04	1.14
1:A:113:SER:HB3	1:A:114:PRO:HD3	1.37	1.06
1:C:1182:GLN:HA	1:C:1192:ARG:HH21	1.22	1.02
1:A:234:ARG:HH11	1:A:234:ARG:HB3	1.21	1.02
1:C:1165:ARG:HD3	2:C:1260:0FP:HA4	1.45	0.97

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	222/256 (87%)	199 (90%)	17 (8%)	6 (3%)	6 16
1	B	222/256 (87%)	200 (90%)	18 (8%)	4 (2%)	11 27
1	C	219/256 (86%)	198 (90%)	17 (8%)	4 (2%)	11 27
1	D	222/256 (87%)	202 (91%)	18 (8%)	2 (1%)	21 49

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	885/1024 (86%)	799 (90%)	70 (8%)	16 (2%)	11	27

5 of 16 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	54	SER
1	A	113	SER
1	A	141	VAL
1	B	413	SER
1	B	441	VAL

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	196/215 (91%)	180 (92%)	16 (8%)	14	32
1	B	195/215 (91%)	178 (91%)	17 (9%)	13	29
1	C	193/215 (90%)	181 (94%)	12 (6%)	23	49
1	D	196/215 (91%)	176 (90%)	20 (10%)	9	21
All	All	780/860 (91%)	715 (92%)	65 (8%)	14	31

5 of 65 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	522	LEU
1	C	1062	ASN
1	D	1515	ARG
1	B	535	LEU
1	C	1041	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 28 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	500	GLN
1	C	1025	GLN
1	D	1443	GLN
1	C	1007	GLN
1	C	1009	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 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	$\# Z > 2$
2	0FP	A	260	1	50,51,51	0.62	1 (2%)	64,71,71	0.82	2 (3%)
2	0FP	B	560	1	50,51,51	0.56	0	64,71,71	0.85	2 (3%)
2	0FP	C	1260	1	50,51,51	0.57	0	64,71,71	0.88	2 (3%)
2	0FP	D	1560	1	50,51,51	0.53	0	64,71,71	0.77	2 (3%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	0FP	A	260	1	-	0/72/72/72	0/1/1/1
2	0FP	B	560	1	-	0/72/72/72	0/1/1/1
2	0FP	C	1260	1	-	0/72/72/72	0/1/1/1
2	0FP	D	1560	1	-	0/72/72/72	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	260	0FP	C4-C11	2.07	1.57	1.52

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1260	0FP	CB2-CA2-C2	-4.12	108.66	112.83
2	A	260	0FP	CB2-CA2-C2	-4.12	108.67	112.83
2	B	560	0FP	CB2-CA2-C2	-3.83	108.95	112.83
2	C	1260	0FP	CB1-CA1-C1	-3.63	109.16	112.83
2	B	560	0FP	CB1-CA1-C1	-3.55	109.24	112.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 30 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	260	0FP	8	0
2	B	560	0FP	6	0
2	C	1260	0FP	14	0
2	D	1560	0FP	2	0

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

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.