wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 09:07 PM GMT

PDB ID	:	1NKK
Title	:	COMPLEX STRUCTURE OF HCMV PROTEASE AND A PEP-TIDOMIMETIC INHIBITOR
Authors	:	Khayat, R.; Batra, R.; Qian, C.; Halmos, T.; Bailey, M.; Tong, L.
Deposited on	:	2003-01-03
Resolution	:	2.60 Å(reported)

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)
Xtriage (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 2.60 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.
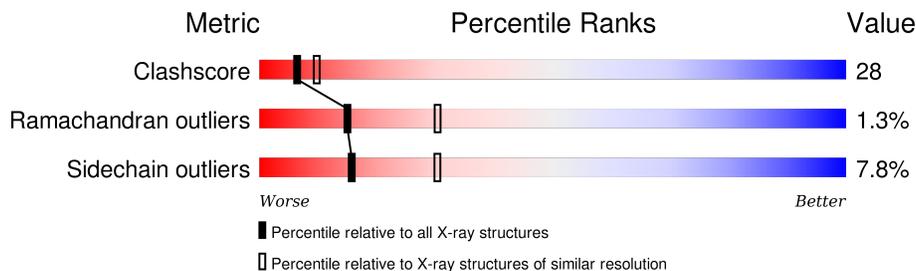
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)

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 <=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	45% 36% 5% 13%
1	B	256	50% 34% • 12%
1	C	256	55% 30% • 12%
1	D	256	48% 33% • 14%
2	E	6	33% 50% 17%
2	F	6	67% 33%
2	G	6	67% 33%

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Mol	Chain	Length	Quality of chain
2	H	6	83% 17%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7431 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 P40.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	222	Total 1749	C 1097	N 324	O 324	S 4	0	0	0
1	B	226	Total 1770	C 1108	N 328	O 328	S 6	0	0	0
1	C	226	Total 1776	C 1112	N 328	O 330	S 6	0	0	0
1	D	220	Total 1731	C 1087	N 319	O 321	S 4	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	31	ARG	GLU	ENGINEERED	UNP P16753
A	143	GLN	ALA	ENGINEERED	UNP P16753
B	331	ARG	GLU	ENGINEERED	UNP P16753
B	443	GLN	ALA	ENGINEERED	UNP P16753
C	1031	ARG	GLU	ENGINEERED	UNP P16753
C	1143	GLN	ALA	ENGINEERED	UNP P16753
D	1331	ARG	GLU	ENGINEERED	UNP P16753
D	1443	GLN	ALA	ENGINEERED	UNP P16753

- Molecule 2 is a protein called Peptidomimetic inhibitor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	6	Total 39	C 23	F 3	N 5	O 8	0	0	0
2	F	6	Total 39	C 23	F 3	N 5	O 8	0	0	0
2	G	6	Total 39	C 23	F 3	N 5	O 8	0	0	0
2	H	6	Total 39	C 23	F 3	N 5	O 8	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	68	Total 68	O 68	0	0
3	B	61	Total 61	O 61	0	0
3	C	64	Total 64	O 64	0	0
3	D	49	Total 49	O 49	0	0
3	E	2	Total 2	O 2	0	0
3	F	3	Total 3	O 3	0	0
3	G	1	Total 1	O 1	0	0
3	H	1	Total 1	O 1	0	0

3 Residue-property plots ⓘ

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

• Molecule 1: Capsid protein P40
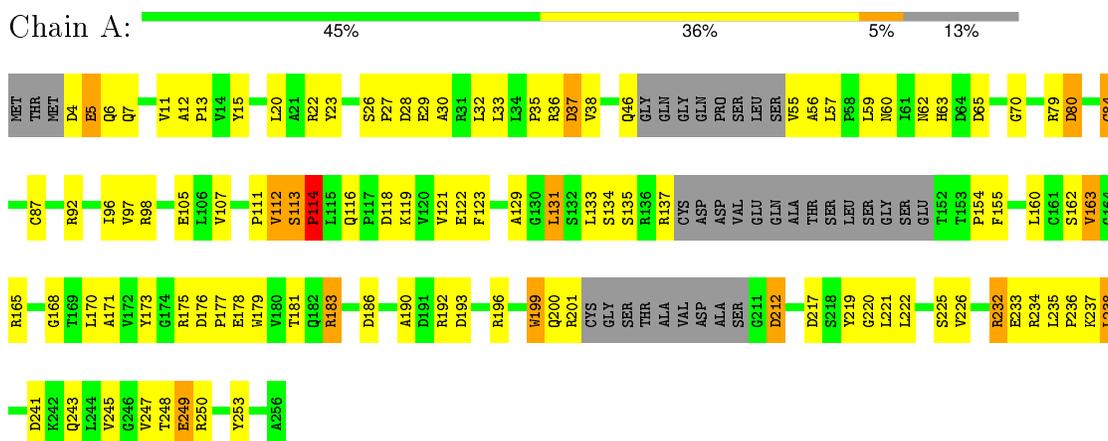
• Molecule 1: Capsid protein P40
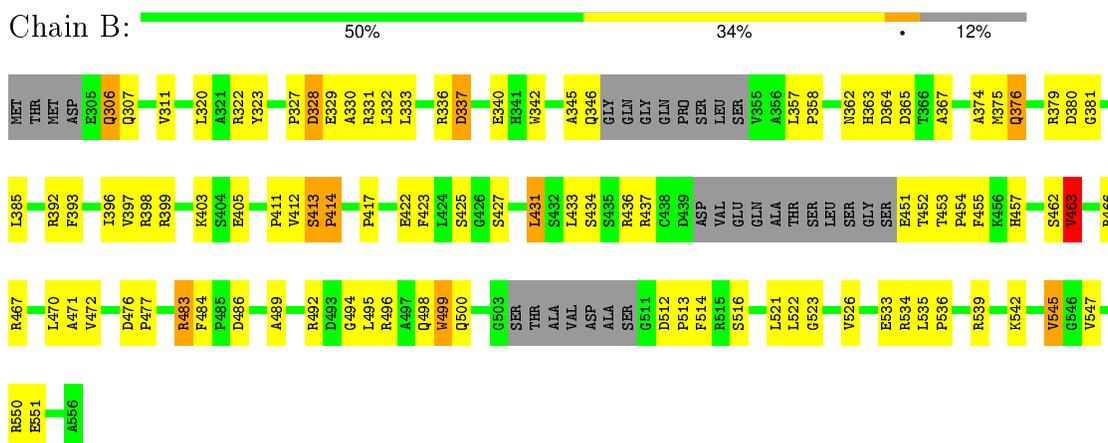
• Molecule 1: Capsid protein P40
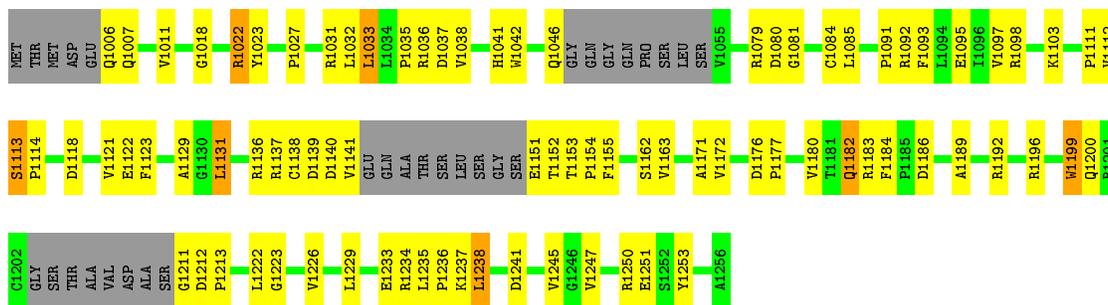
● Molecule 1: Capsid protein P40

Chain D:
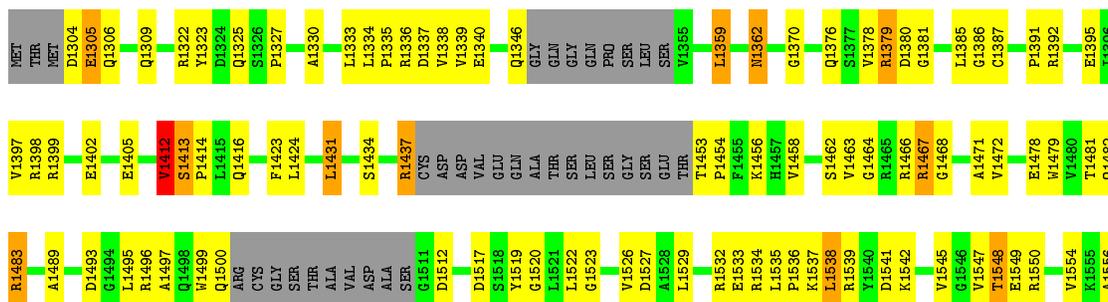
● Molecule 2: Peptidomimetic inhibitor

Chain E:

● Molecule 2: Peptidomimetic inhibitor

Chain F:

● Molecule 2: Peptidomimetic inhibitor

Chain G:

● Molecule 2: Peptidomimetic inhibitor

Chain H:

4 Data and refinement statistics (i)

Xtriage (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	105.17Å 215.25Å 52.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.98 – 2.60	Depositor
% Data completeness (in resolution range)	82.6 (19.98-2.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.233 , 0.280	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7431	wwPDB-VP
Average B, all atoms (Å^2)	17.0	wwPDB-VP

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: DMH, CFT, DMK, ACE

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	0.39	0/1784	0.62	0/2417
1	B	0.41	0/1805	0.63	0/2445
1	C	0.38	0/1811	0.62	0/2454
1	D	0.38	0/1766	0.61	0/2393
2	E	2.53	1/11 (9.1%)	1.68	1/13 (7.7%)
2	F	2.78	1/11 (9.1%)	1.62	1/13 (7.7%)
2	G	2.56	1/11 (9.1%)	1.52	1/13 (7.7%)
2	H	2.61	1/11 (9.1%)	1.75	1/13 (7.7%)
All	All	0.44	4/7210 (0.1%)	0.63	4/9761 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	564	ALA	C-O	8.90	1.40	1.23
2	H	1564	ALA	C-O	8.57	1.39	1.23
2	E	264	ALA	C-O	8.17	1.38	1.23
2	G	1264	ALA	C-O	8.17	1.38	1.23

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	E	264	ALA	CA-C-O	-5.77	107.97	120.10
2	H	1564	ALA	CA-C-O	-5.71	108.10	120.10
2	F	564	ALA	CA-C-O	-5.40	108.76	120.10
2	G	1264	ALA	CA-C-O	-5.19	109.20	120.10

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	1749	0	1731	108	0
1	B	1770	0	1746	107	0
1	C	1776	0	1754	85	0
1	D	1731	0	1711	101	0
2	E	39	0	33	4	0
2	F	39	0	33	1	0
2	G	39	0	33	1	0
2	H	39	0	33	0	0
3	A	68	0	0	6	0
3	B	61	0	0	9	0
3	C	64	0	0	2	0
3	D	49	0	0	5	0
3	E	2	0	0	1	0
3	F	3	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
All	All	7431	0	7074	393	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 28.

The worst 5 of 393 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:C:1182:GLN:HE21	1:C:1182:GLN:HA	1.16	1.09
1:D:1413:SER:HB3	1:D:1414:PRO:HD3	1.37	1.07
1:C:1042:TRP:HE1	1:C:1140:ASP:HB3	1.11	1.06
1:B:462:SER:HB2	1:B:534:ARG:HE	1.17	1.03
1:B:331:ARG:HD3	1:B:437:ARG:HH21	1.22	1.03

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	214/256 (84%)	199 (93%)	11 (5%)	4 (2%)	10	19
1	B	218/256 (85%)	198 (91%)	18 (8%)	2 (1%)	21	42
1	C	218/256 (85%)	198 (91%)	17 (8%)	3 (1%)	14	28
1	D	212/256 (83%)	190 (90%)	20 (9%)	2 (1%)	21	42
2	E	1/6 (17%)	1 (100%)	0	0	100	100
2	F	1/6 (17%)	1 (100%)	0	0	100	100
2	G	1/6 (17%)	1 (100%)	0	0	100	100
2	H	1/6 (17%)	1 (100%)	0	0	100	100
All	All	866/1048 (83%)	789 (91%)	66 (8%)	11 (1%)	15	30

5 of 11 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	113	SER
1	B	413	SER
1	C	1113	SER
1	D	1413	SER
1	A	212	ASP

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	188/215 (87%)	170 (90%)	18 (10%)	10	20

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	190/215 (88%)	175 (92%)	15 (8%)	15	30
1	C	192/215 (89%)	183 (95%)	9 (5%)	32	59
1	D	186/215 (86%)	169 (91%)	17 (9%)	12	22
2	E	1/1 (100%)	1 (100%)	0	100	100
2	F	1/1 (100%)	1 (100%)	0	100	100
2	G	1/1 (100%)	1 (100%)	0	100	100
2	H	1/1 (100%)	1 (100%)	0	100	100
All	All	760/864 (88%)	701 (92%)	59 (8%)	16	30

5 of 59 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	414	PRO
1	C	1022	ARG
1	D	1467	ARG
1	B	431	LEU
1	B	483	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	482	GLN
1	C	1007	GLN
1	D	1362	ASN
1	B	325	GLN
1	B	346	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)

8 non-standard protein/DNA/RNA residues 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	DMK	E	262	2	5,9,10	1.53	1 (20%)	5,13,15	0.78	0
2	DMH	E	263	2	8,9,10	0.58	0	7,11,13	0.86	0
2	DMK	F	562	2	5,9,10	1.06	1 (20%)	5,13,15	1.11	0
2	DMH	F	563	2	8,9,10	0.73	0	7,11,13	0.85	0
2	DMK	G	1262	2	5,9,10	1.34	1 (20%)	5,13,15	0.80	0
2	DMH	G	1263	2	8,9,10	0.68	0	7,11,13	0.85	0
2	DMK	H	1562	2	5,9,10	1.44	1 (20%)	5,13,15	0.99	1 (20%)
2	DMH	H	1563	2	8,9,10	0.69	0	7,11,13	0.81	0

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	DMK	E	262	2	-	0/4/14/16	0/0/0/0
2	DMH	E	263	2	-	0/8/10/12	0/0/0/0
2	DMK	F	562	2	-	0/4/14/16	0/0/0/0
2	DMH	F	563	2	-	0/8/10/12	0/0/0/0
2	DMK	G	1262	2	-	0/4/14/16	0/0/0/0
2	DMH	G	1263	2	-	0/8/10/12	0/0/0/0
2	DMK	H	1562	2	-	0/4/14/16	0/0/0/0
2	DMH	H	1563	2	-	0/8/10/12	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	562	DMK	CB-CG1	2.00	1.54	1.50
2	G	1262	DMK	CB-CG1	2.61	1.55	1.50
2	H	1562	DMK	CB-CG1	2.68	1.55	1.50
2	E	262	DMK	CB-CG1	2.93	1.56	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	1562	DMK	CG3-CB-CG2	2.10	109.91	107.33

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	262	DMK	1	0
2	F	563	DMH	1	0
2	G	1262	DMK	1	0

5.5 Carbohydrates (i)

There are no carbohydrates 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)

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

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

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

6.3 Carbohydrates (i)

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

6.4 Ligands (i)

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

6.5 Other polymers (i)

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