



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 04:49 PM GMT

PDB ID : 4G93
Title : CRYSTAL STRUCTURE OF THE HUMAN HEPATITIS B VIRUS T = 4
CAPSID, ADYW STRAIN, in COMPLEX WITH THE PHENYLPROPE-
NAMIDE ASSEMBLY ACCELERATOR AT-130
Authors : Katen, S.P.; Zlotnick, A.
Deposited on : 2012-07-23
Resolution : 4.20 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)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

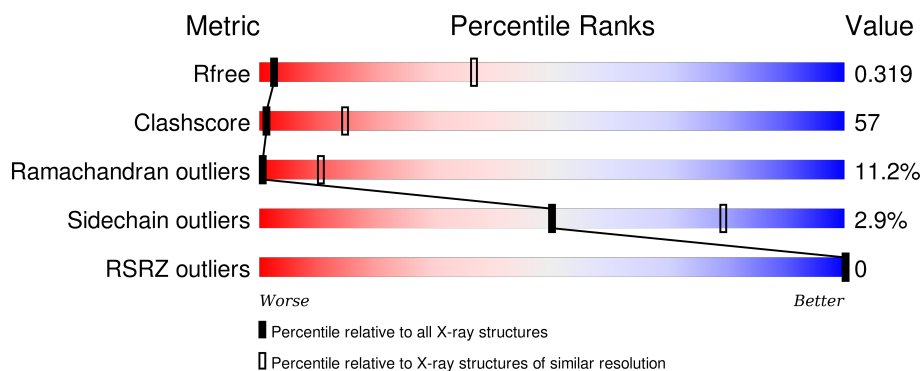
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.20 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}	91344	1039 (4.80-3.60)
Clashscore	102246	1140 (4.80-3.60)
Ramachandran outliers	100387	1083 (4.80-3.60)
Sidechain outliers	100360	1067 (4.80-3.60)
RSRZ outliers	91569	1042 (4.80-3.60)

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 $\leq 5\%$. 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	150	<div> <div>34%</div> <div>53%</div> <div>7%</div> <div>5%</div> </div>
1	B	150	<div> <div>30%</div> <div>51%</div> <div>12%</div> <div>7%</div> </div>
1	C	150	<div> <div>31%</div> <div>55%</div> <div>6%</div> <div>7%</div> </div>
1	D	150	<div> <div>28%</div> <div>54%</div> <div>13%</div> <div>5%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	0YP	B	201	-	-	X	X
2	0YP	C	201	-	-	-	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4544 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.

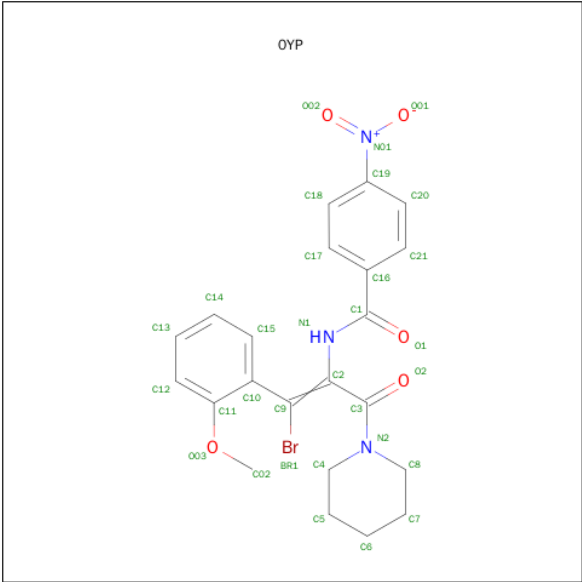
- Molecule 1 is a protein called Capsid protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	142	Total	C	N	O	S	0	0	0
			1129	735	186	206	2			
1	B	140	Total	C	N	O	S	0	0	0
			1116	728	184	202	2			
1	C	139	Total	C	N	O	S	0	0	0
			1108	722	183	201	2			
1	D	142	Total	C	N	O	S	0	0	0
			1129	735	186	206	2			

There are 16 discrepancies between the modelled and reference sequences:

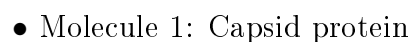
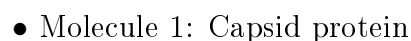
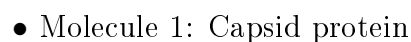
Chain	Residue	Modelled	Actual	Comment	Reference
A	48	ALA	CYS	ENGINEERED MUTATION	UNP P03147
A	61	ALA	CYS	ENGINEERED MUTATION	UNP P03147
A	107	ALA	CYS	ENGINEERED MUTATION	UNP P03147
A	150	CYS	-	SEE REMARK 999	UNP P03147
B	48	ALA	CYS	ENGINEERED MUTATION	UNP P03147
B	61	ALA	CYS	ENGINEERED MUTATION	UNP P03147
B	107	ALA	CYS	ENGINEERED MUTATION	UNP P03147
B	150	CYS	-	SEE REMARK 999	UNP P03147
C	48	ALA	CYS	ENGINEERED MUTATION	UNP P03147
C	61	ALA	CYS	ENGINEERED MUTATION	UNP P03147
C	107	ALA	CYS	ENGINEERED MUTATION	UNP P03147
C	150	CYS	-	SEE REMARK 999	UNP P03147
D	48	ALA	CYS	ENGINEERED MUTATION	UNP P03147
D	61	ALA	CYS	ENGINEERED MUTATION	UNP P03147
D	107	ALA	CYS	ENGINEERED MUTATION	UNP P03147
D	150	CYS	-	SEE REMARK 999	UNP P03147

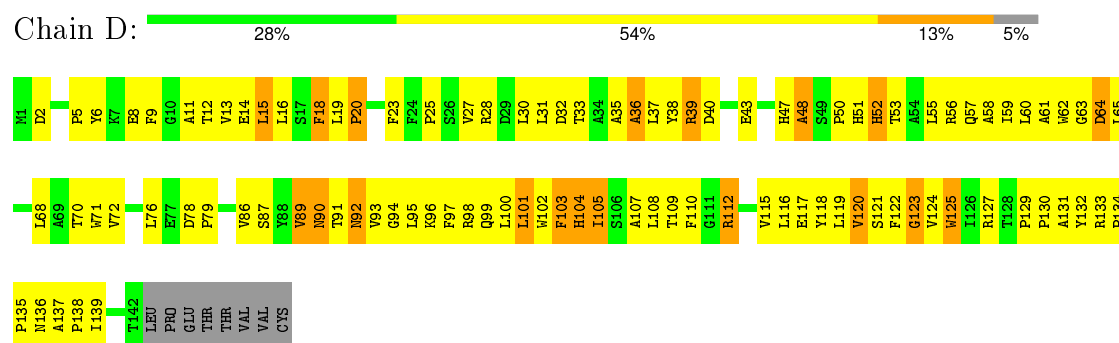
- Molecule 2 is N-[(1E)-1-BROMO-1-(2-METHOXYPHENYL)-3-OXO-3-(PIPERIDIN-1-YL)PROP-1-EN-2-YL]-4-NITROBENZAMIDE (three-letter code: 0YP) (formula: C₂₂H₂₂BrN₃O₅).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	Br	C	N	O	5	0
			31	1	22	3	5		
2	C	1	Total	Br	C	N	O	5	0
			31	1	22	3	5		

- Molecule 1: Capsid protein





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	527.38Å 362.75Å 538.23Å 90.00° 105.13° 90.00°	Depositor
Resolution (Å)	40.00 – 4.20 29.99 – 3.91	Depositor EDS
% Data completeness (in resolution range)	71.1 (40.00-4.20) 72.7 (29.99-3.91)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.04 (at 3.86Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.379 , 0.383 0.327 , 0.319	Depositor DCC
R_{free} test set	5100 reflections (0.98%)	DCC
Wilson B-factor (Å ²)	141.3	Xtriage
Anisotropy	0.149	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.16 , 6.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	3 of 635269 reflections (0.000%)	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4544	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.58% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 0YP

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	0.33	0/1164	0.55	1/1597 (0.1%)
1	B	0.32	0/1151	0.54	0/1579
1	C	0.36	0/1143	0.51	0/1568
1	D	0.32	0/1164	0.53	1/1597 (0.1%)
All	All	0.33	0/4622	0.53	2/6341 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	64	ASP	CB-CG-OD2	5.21	122.99	118.30
1	A	64	ASP	CB-CG-OD2	5.13	122.92	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	A	1129	0	1110	121	0
1	B	1116	0	1098	149	0
1	C	1108	0	1087	138	0
1	D	1129	0	1110	124	0
2	B	31	0	22	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	31	0	22	4	0
All	All	4544	0	4449	512	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 57.

The worst 5 of 512 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:30:LEU:HD21	1:C:101:LEU:HB3	1.20	1.18
1:C:30:LEU:HD21	1:C:101:LEU:CB	1.78	1.12
1:B:126:ILE:HG23	1:B:127:ARG:HD3	1.41	1.03
1:B:13:VAL:HA	1:B:16:LEU:HD23	1.39	1.02
2:B:201:OYP:H21	1:C:128:THR:HG22	1.43	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	140/150 (93%)	86 (61%)	41 (29%)	13 (9%)	1	16
1	B	138/150 (92%)	78 (56%)	43 (31%)	17 (12%)	0	8
1	C	137/150 (91%)	83 (61%)	42 (31%)	12 (9%)	1	17
1	D	140/150 (93%)	73 (52%)	47 (34%)	20 (14%)	0	6
All	All	555/600 (92%)	320 (58%)	173 (31%)	62 (11%)	0	10

5 of 62 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	89	VAL

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Mol	Chain	Res	Type
1	A	101	LEU
1	A	103	PHE
1	A	105	ILE
1	B	7	LYS

5.3.2 Protein sidechains ⓘ

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	122/130 (94%)	121 (99%)	1 (1%)	86	93
1	B	120/130 (92%)	116 (97%)	4 (3%)	45	78
1	C	119/130 (92%)	115 (97%)	4 (3%)	44	77
1	D	122/130 (94%)	117 (96%)	5 (4%)	37	73
All	All	483/520 (93%)	469 (97%)	14 (3%)	50	79

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	9	PHE
1	C	18	PHE
1	D	40	ASP
1	C	2	ASP
1	D	18	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	47	HIS
1	C	75	ASN
1	D	52	HIS
1	B	99	GLN
1	C	136	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

2 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	0YP	B	201	-	29,33,33	3.20	11 (37%)	36,45,45	2.37	8 (22%)
2	0YP	C	201	-	29,33,33	3.28	11 (37%)	36,45,45	1.89	6 (16%)

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	0YP	B	201	-	-	0/24/38/38	0/3/3/3
2	0YP	C	201	-	-	2/24/38/38	0/3/3/3

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	201	0YP	C8-N2	-4.99	1.38	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	201	0YP	C8-N2	-4.35	1.39	1.47
2	B	201	0YP	C4-N2	-3.02	1.41	1.47
2	C	201	0YP	C4-N2	-2.49	1.42	1.47
2	B	201	0YP	C1-N1	3.55	1.44	1.37

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	201	0YP	O03-C11-C12	-5.83	114.53	124.35
2	C	201	0YP	O2-C3-N2	-5.55	113.26	122.50
2	B	201	0YP	C2-N1-C1	-4.60	109.81	122.72
2	B	201	0YP	O2-C3-N2	-4.56	114.91	122.50
2	C	201	0YP	O03-C11-C12	-4.54	116.70	124.35

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	201	0YP	C20-C19-N01-O02
2	C	201	0YP	C18-C19-N01-O02

There are no ring outliers.

2 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	201	0YP	10	0
2	C	201	0YP	4	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 ⓘ

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	142/150 (94%)	-0.88	0 100 100	35, 80, 150, 166	0
1	B	140/150 (93%)	-0.88	0 100 100	34, 77, 163, 177	0
1	C	139/150 (92%)	-0.78	0 100 100	29, 82, 208, 224	0
1	D	142/150 (94%)	-0.79	0 100 100	35, 77, 195, 208	0
All	All	563/600 (93%)	-0.83	0 100 100	29, 79, 173, 224	0

There are no RSRZ outliers to report.

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	0YP	C	201	31/31	0.86	0.45	3.83	0,150,152,153	5
2	0YP	B	201	31/31	0.95	0.30	3.21	0,62,65,70	5

6.5 Other polymers [i](#)

There are no such residues in this entry.