



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 03:52 pm GMT

PDB ID : 1O55
Title : Crystal structure of HCV NS5B RNA polymerase complexed with a novel non-competitive inhibitor.
Authors : Love, R.A.; Parge, H.E.; Yu, X.; Hickey, M.J.; Diehl, W.; Gao, J.; Wriggers, H.; Ekker, A.; Wang, L.; Thomson, J.A.; Dragovich, P.S.; Fuhrman, S.A.
Deposited on : 2003-03-18
Resolution : 2.20 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

<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 : 1.7.2 (RC1), CSD as538be (2017)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

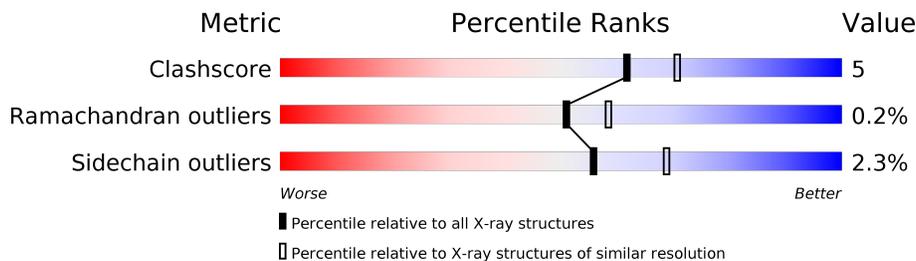
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)

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	576	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5009 atoms, of which 37 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.

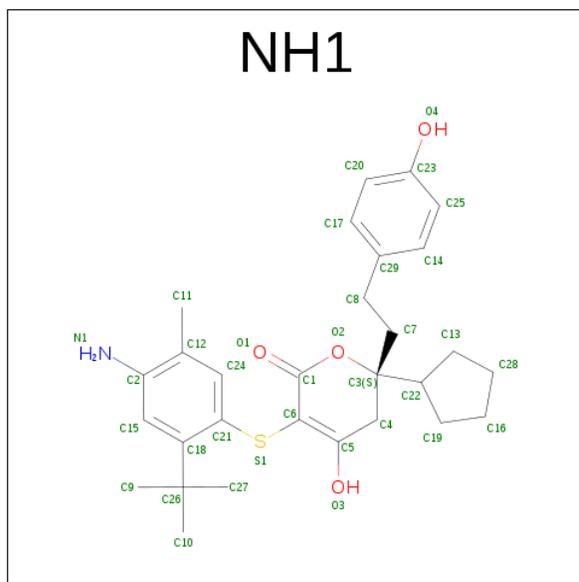
- Molecule 1 is a protein called Hepatitis C virus NS5B RNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	563	4377	2757	774	815	31	0	0	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	GLN	LEU	ENGINEERED	UNP P26663
A	101	TYR	PHE	ENGINEERED	UNP P26663
A	114	ARG	LYS	ENGINEERED	UNP P26663
A	572	HIS	-	EXPRESSION TAG	UNP P26663
A	573	HIS	-	EXPRESSION TAG	UNP P26663
A	574	HIS	-	EXPRESSION TAG	UNP P26663
A	575	HIS	-	EXPRESSION TAG	UNP P26663
A	576	HIS	-	EXPRESSION TAG	UNP P26663
A	577	HIS	-	EXPRESSION TAG	UNP P26663

- Molecule 2 is 3-(4-AMINO-2-TERT-BUTYL-5-METHYL-PHENYLSULFANYL)-6-CYCLOPENTYL-4-HYDROXY-6-[2-(4-HYDROXY-PHENYL)-ETHYL]-5,6-DIHYDRO-PYRAZIN-2-ONE (three-letter code: NH1) (formula: C₂₉H₃₇NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	H	N	O			S
2	A	1	72	29	37	1	4	1	0	0

- Molecule 3 is water.

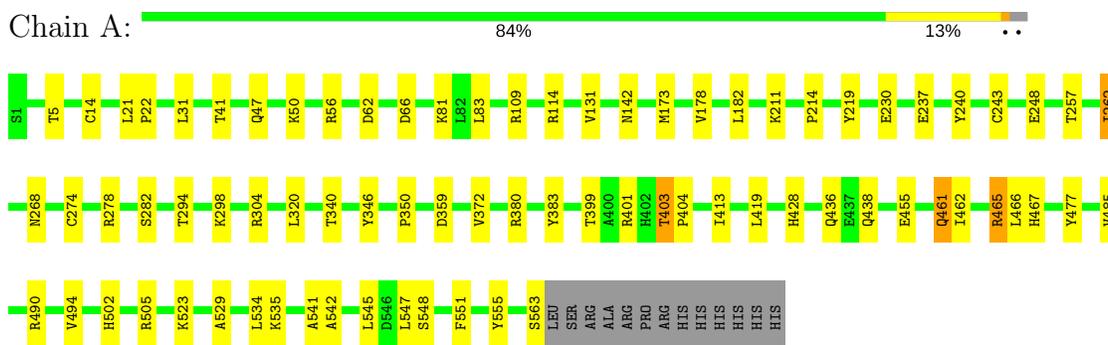
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	560	Total	O	0	0
			560	560		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

Note EDS was not executed.

- Molecule 1: Hepatitis C virus NS5B RNA polymerase



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	83.00Å 83.00Å 180.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.20	Depositor
% Data completeness (in resolution range)	97.0 (10.00-2.20)	Depositor
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.207 , 0.251	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	5009	wwPDB-VP
Average B, all atoms (Å ²)	18.0	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: NH1

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	0.68	0/4473	0.87	5/6071 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	240	TYR	CB-CG-CD2	-7.53	116.48	121.00
1	A	346	TYR	CB-CG-CD2	-6.20	117.28	121.00
1	A	535	LYS	CA-CB-CG	-5.75	100.76	113.40
1	A	555	TYR	CB-CG-CD2	-5.45	117.73	121.00
1	A	304	ARG	NE-CZ-NH2	-5.07	117.77	120.30

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	4377	0	4389	43	1
2	A	35	37	36	2	0
3	A	560	0	0	11	0
All	All	4972	37	4425	45	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 5.

All (45) 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:436:GLN:HB2	1:A:438:GLN:HG3	1.66	0.78
1:A:219:TYR:HB3	1:A:320:LEU:HD23	1.79	0.64
1:A:182:LEU:HD12	1:A:243:CYS:SG	2.37	0.64
1:A:83:LEU:HB2	1:A:173:MET:HA	1.83	0.61
1:A:413:ILE:O	1:A:467:HIS:HE1	1.82	0.61
1:A:5:THR:HG23	1:A:278:ARG:HH12	1.67	0.60
1:A:178:VAL:HG23	3:A:1216:HOH:O	2.03	0.59
1:A:14:CYS:HB2	3:A:1108:HOH:O	2.03	0.58
1:A:383:TYR:HH	1:A:477:TYR:HD2	1.51	0.58
1:A:545:LEU:HB3	1:A:547:LEU:HD13	1.84	0.58
1:A:268:ASN:HB3	1:A:274:CYS:SG	2.43	0.57
1:A:461:GLN:HG2	1:A:541:ALA:HB3	1.85	0.57
1:A:455:GLU:HB3	3:A:1248:HOH:O	2.05	0.57
1:A:230:GLU:HG3	1:A:262:ILE:HG13	1.87	0.55
2:A:901:NH1:S1	2:A:901:NH1:H23	2.47	0.55
1:A:428:HIS:HD2	3:A:1021:HOH:O	1.91	0.53
1:A:462:ILE:O	1:A:466:LEU:HB2	2.09	0.52
1:A:399:THR:OG1	1:A:428:HIS:HE1	1.93	0.52
1:A:523:LYS:HG3	1:A:534:LEU:HD12	1.91	0.51
1:A:502:HIS:HE1	3:A:1217:HOH:O	1.93	0.51
1:A:340:THR:CG2	1:A:350:PRO:HG3	2.41	0.51
1:A:83:LEU:CB	1:A:173:MET:HA	2.40	0.51
1:A:465:ARG:HH12	1:A:545:LEU:N	2.09	0.51
1:A:419:LEU:HD11	1:A:485:VAL:HG11	1.94	0.49
1:A:237:GLU:HG3	1:A:257:THR:OG1	2.13	0.49
1:A:109:ARG:HH11	1:A:109:ARG:HG2	1.77	0.48
1:A:66:ASP:HB3	3:A:1292:HOH:O	2.13	0.48
1:A:50:LYS:HE3	3:A:1402:HOH:O	2.14	0.48
1:A:31:LEU:HB3	1:A:494:VAL:HG22	1.96	0.47
1:A:47:GLN:HG3	3:A:1375:HOH:O	2.14	0.47
2:A:901:NH1:S1	2:A:901:NH1:H16	2.55	0.47
1:A:294:THR:HG23	1:A:298:LYS:HE3	1.98	0.46
1:A:372:VAL:CG2	1:A:380:ARG:HG3	2.46	0.46
1:A:403:THR:HA	1:A:404:PRO:HD2	1.79	0.46
1:A:21:LEU:HD12	1:A:22:PRO:HD2	1.97	0.46
1:A:211:LYS:HB2	1:A:214:PRO:HB3	1.98	0.46
1:A:41:THR:HA	1:A:142:ASN:OD1	2.19	0.43
1:A:466:LEU:HD11	1:A:551:PHE:HE2	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:372:VAL:HG23	1:A:380:ARG:HG3	2.00	0.42
1:A:461:GLN:HB3	1:A:542:ALA:HA	2.01	0.42
1:A:62:ASP:HB2	3:A:1453:HOH:O	2.20	0.41
1:A:248:GLU:HB2	3:A:1279:HOH:O	2.19	0.41
1:A:467:HIS:HD2	3:A:910:HOH:O	2.03	0.41
1:A:505:ARG:HG3	1:A:529:ALA:HB1	2.02	0.40
1:A:257:THR:HG22	1:A:262:ILE:HD13	2.03	0.40

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:359:ASP:OD2	1:A:401:ARG:NH2[6_555]	2.17	0.03

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	561/576 (97%)	545 (97%)	15 (3%)	1 (0%)	51 58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	131	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	479/491 (98%)	468 (98%)	11 (2%)	56 69

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	ARG
1	A	81	LYS
1	A	114	ARG
1	A	262	ILE
1	A	282	SER
1	A	403	THR
1	A	461	GLN
1	A	465	ARG
1	A	490	ARG
1	A	548	SER
1	A	563	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	ASN
1	A	428	HIS
1	A	446	GLN
1	A	467	HIS
1	A	514	GLN
1	A	544	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

1 ligand is 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	NH1	A	901	-	35,38,38	1.09	4 (11%)	46,57,57	1.23	3 (6%)

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	NH1	A	901	-	-	0/22/47/47	0/4/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	NH1	C26-C18	-2.79	1.49	1.54
2	A	901	NH1	C11-C12	-2.12	1.46	1.51
2	A	901	NH1	C21-C18	2.08	1.42	1.40
2	A	901	NH1	C7-C3	2.41	1.56	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	NH1	C3-C4-C5	-3.28	105.73	113.35
2	A	901	NH1	C13-C22-C3	2.24	115.87	112.31
2	A	901	NH1	O2-C1-O1	4.63	123.48	117.69

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	NH1	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.