



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 06:28 PM GMT

PDB ID : 4LQ3  
Title : Crystal structure of human norovirus RNA-dependent RNA-polymerase bound to the inhibitor PPNDs  
Authors : Milani, M.; Tarantino, D.; Mastrangelo, E.; Croci, R.  
Deposited on : 2013-07-17  
Resolution : 2.60 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](mailto: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.

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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

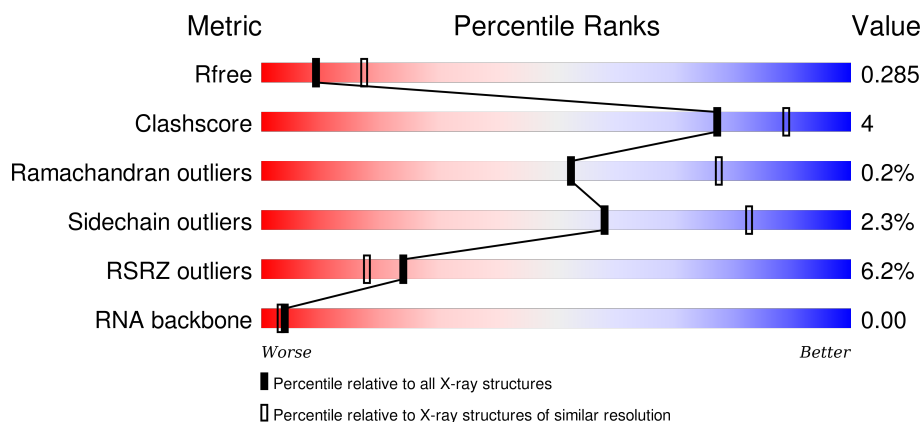
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


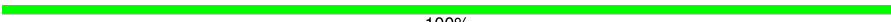
The reported resolution of this entry is 2.60 Å.

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))
$R_{free}$	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)
RNA backbone	2183	1022 (3.00-2.20)

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  $\geq 3$ , 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	526	 6% 86% 7% • 6%
2	R	2	 100%

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4215 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 RNA-dependent RNA-polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	495	Total	C	N	O	S	0	0	0
			3880	2463	664	731	22			

There are 19 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	EXPRESSION TAG	UNP Q2N379
A	1	GLY	-	EXPRESSION TAG	UNP Q2N379
A	2	GLY	-	EXPRESSION TAG	UNP Q2N379
A	153	GLY	GLU	CONFLICT	UNP Q2N379
A	511	VAL	-	EXPRESSION TAG	UNP Q2N379
A	512	ASP	-	EXPRESSION TAG	UNP Q2N379
A	513	LYS	-	EXPRESSION TAG	UNP Q2N379
A	514	LEU	-	EXPRESSION TAG	UNP Q2N379
A	515	ALA	-	EXPRESSION TAG	UNP Q2N379
A	516	ALA	-	EXPRESSION TAG	UNP Q2N379
A	517	ALA	-	EXPRESSION TAG	UNP Q2N379
A	518	LEU	-	EXPRESSION TAG	UNP Q2N379
A	519	GLU	-	EXPRESSION TAG	UNP Q2N379
A	520	HIS	-	EXPRESSION TAG	UNP Q2N379
A	521	HIS	-	EXPRESSION TAG	UNP Q2N379
A	522	HIS	-	EXPRESSION TAG	UNP Q2N379
A	523	HIS	-	EXPRESSION TAG	UNP Q2N379
A	524	HIS	-	EXPRESSION TAG	UNP Q2N379
A	525	HIS	-	EXPRESSION TAG	UNP Q2N379

- Molecule 2 is a RNA chain called 5'-R(P\*GP\*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	R	2	Total	C	N	O	P	0	0	0
			46	20	10	14	2			

- 20V
- 
- The image displays a complex organic molecule, likely a derivative of a heterocyclic compound, featuring a central benzene ring substituted with various functional groups and labeled atoms. The structure includes a nitro group (NO<sub>2</sub>) at the top left, a sulfonamide group (SO<sub>2</sub>NH<sub>2</sub>) at the top right, a sulfonamide group (SO<sub>2</sub>NH<sub>2</sub>) at the bottom left, and a pyridine ring at the bottom right. The molecule is heavily labeled with abbreviations such as OAI, OAJ, NBJ, CBB, CAQ, CAG, SBM, OAN, OAF, CBI, CBH, CAS, CBA, CAY, CAA, CBF, CAP, CAZ, NAV, NAW, CBD, CBE, CBC, CAO, OAB, OAK, OAL, OAE, OAC, PBK, CAT, and OAX. The labels are color-coded: blue for nitrogen-containing groups, green for sulfur-containing groups, red for oxygen-containing groups, and black for carbon-containing groups. The molecule is shown in a 2D representation with a white background.

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

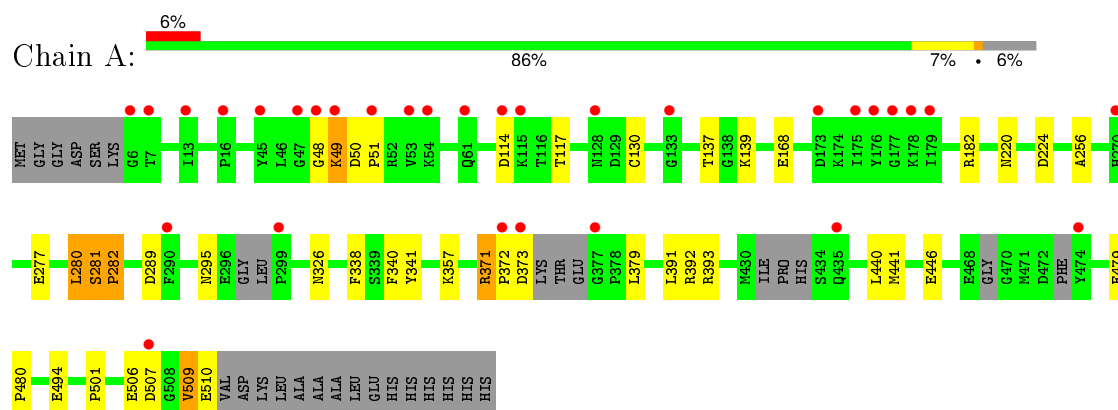
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	129	Total O 129 129	0	0
5	R	3	Total O 3 3	0	0

### 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 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.

#### • Molecule 1: RNA-dependent RNA-polymerase



#### • Molecule 2: 5'-R(P\*GP\*G)-3'



There are no outlier residues recorded for this chain.

## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.75Å 116.10Å 122.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.40 – 2.60 38.41 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.6 (38.40-2.60) 99.7 (38.41-2.60)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.79 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.220 , 0.285 0.222 , 0.285	Depositor DCC
$R_{free}$ test set	985 reflections (5.45%)	DCC
Wilson B-factor (Å <sup>2</sup> )	58.0	Xtriage
Anisotropy	0.182	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 44.1	EDS
Estimated twinning fraction	0.015 for -h,-l,-k	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 19046 reflections (0.005%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4215	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.76% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MG, 20V

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	0.41	2/3974 (0.1%)	0.57	2/5381 (0.0%)
2	R	0.22	0/51	0.61	0/78
All	All	0.41	2/4025 (0.0%)	0.57	2/5459 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	51	PRO	N-CD	5.22	1.55	1.47
1	A	282	PRO	N-CD	5.13	1.55	1.47

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	50	ASP	C-N-CD	5.67	140.31	128.40
1	A	281	SER	C-N-CD	5.64	140.25	128.40

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	3880	0	3850	28	0
2	R	46	0	23	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	156	0	57	6	0
4	A	1	0	0	0	0
5	A	129	0	0	0	0
5	R	3	0	0	0	0
All	All	4215	0	3930	30	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 4.

The worst 5 of 30 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:509:VAL:HA	1:A:510:GLU:C	2.09	0.72
1:A:393:ARG:NH2	1:A:446:GLU:OE1	2.30	0.64
3:A:602:20V:OAM	3:A:602:20V:H12	1.98	0.63
3:A:602:20V:OAG	3:A:602:20V:H10	2.01	0.60
1:A:256:ALA:HA	1:A:280:LEU:HD21	1.84	0.60

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	483/526 (92%)	475 (98%)	7 (1%)	1 (0%)	52 77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	371	ARG



### 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	427/451 (95%)	417 (98%)	10 (2%)	58 83

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	326	ASN
1	A	341	TYR
1	A	440	LEU
1	A	289	ASP
1	A	373	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	96	GLN
1	A	143	GLN
1	A	148	ASN
1	A	326	ASN
1	A	423	HIS

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	R	1/2 (50%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

## 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

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

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$
3	20V	A	601	-	39,41,41	2.70	14 (35%)	48,64,64	1.76	13 (27%)
3	20V	A	602	-	39,41,41	2.62	14 (35%)	48,64,64	1.97	9 (18%)
3	20V	A	603	-	39,41,41	2.88	14 (35%)	48,64,64	1.66	13 (27%)
3	20V	A	604	-	39,41,41	3.19	13 (33%)	48,64,64	1.58	10 (20%)

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
3	20V	A	601	-	-	0/28/29/29	0/3/3/3
3	20V	A	602	-	-	0/28/29/29	0/3/3/3
3	20V	A	603	-	-	0/28/29/29	0/3/3/3
3	20V	A	604	-	-	2/28/29/29	0/3/3/3

The worst 5 of 55 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	603	20V	CBD-NAV	-5.87	1.33	1.41
3	A	602	20V	CAT-CBE	-5.59	1.40	1.51
3	A	601	20V	CAA-CAY	-5.43	1.39	1.50
3	A	604	20V	CBD-NAV	-5.37	1.34	1.41
3	A	602	20V	CBD-NAV	-5.33	1.34	1.41

The worst 5 of 45 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	20V	OAB-CAO-CBC	-3.07	118.90	125.11
3	A	604	20V	CAS-CBI-CBF	-2.70	120.12	123.23
3	A	601	20V	CBA-CBC-CBE	-2.68	117.65	120.02
3	A	601	20V	CAS-CBI-CBF	-2.58	120.25	123.23
3	A	602	20V	CBA-CAY-NAW	-2.58	117.48	120.71

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	604	20V	CAQ-CBB-NBJ-OAI
3	A	604	20V	CAS-CBB-NBJ-OAI

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	20V	4	0
3	A	603	20V	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](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	495/526 (94%)	0.14	31 (6%) 23 17	26, 53, 82, 107	5 (1%)
2	R	2/2 (100%)	1.26	0 100 100	93, 93, 93, 106	0
All	All	497/528 (94%)	0.15	31 (6%) 24 18	26, 53, 83, 107	5 (1%)

The worst 5 of 31 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	6	GLY	5.6
1	A	54	LYS	4.1
1	A	48	GLY	4.1
1	A	474	TYR	4.1
1	A	7	THR	3.6

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

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

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
3	20V	A	603	39/39	0.87	0.21	0.55	70,82,89,92	0
4	MG	A	605	1/1	0.77	0.20	0.48	45,45,45,45	0
3	20V	A	604	39/39	0.86	0.22	-0.13	91,101,111,112	0
3	20V	A	602	39/39	0.93	0.14	-0.44	49,59,70,75	0
3	20V	A	601	39/39	0.95	0.13	-1.38	42,53,72,81	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.