



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 12:52 am GMT

PDB ID : 3V7A
Title : Structural basis for broad detection of genogroup II noroviruses by a monoclonal antibody that binds to a site occluded in the viral particle
Authors : Hansman, G.S.; Mclellan, J.S.; Kwong, P.D.
Deposited on : 2011-12-20
Resolution : 3.30 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
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

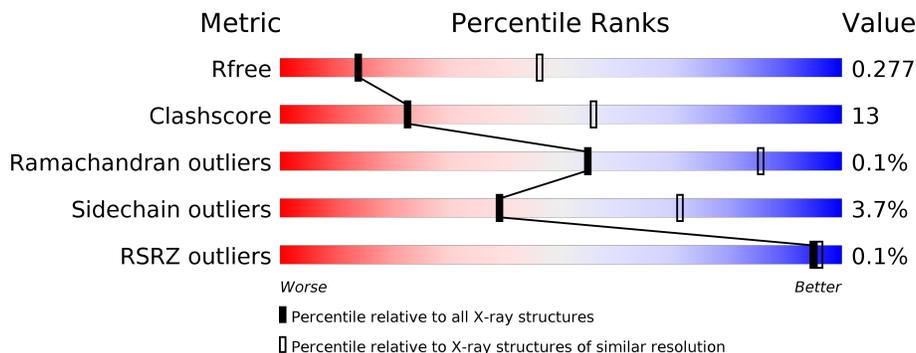
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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}	100719	1034 (3.36-3.24)
Clashscore	112137	1100 (3.36-3.24)
Ramachandran outliers	110173	1081 (3.36-3.24)
Sidechain outliers	110143	1080 (3.36-3.24)
RSRZ outliers	101464	1039 (3.36-3.24)

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	E	223	 65% 31%
1	F	223	 65% 31%
2	G	215	 72% 27%
2	H	215	 74% 25%
3	A	315	 71% 28%
3	B	315	 72% 25%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11459 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 5B18 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	F	220	Total 1643	C 1035	N 273	O 328	S 7	0	0	0
1	E	220	Total 1643	C 1035	N 273	O 328	S 7	0	0	0

- Molecule 2 is a protein called 5B18 kappa chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	G	215	Total 1671	C 1044	N 281	O 339	S 7	0	0	0
2	H	215	Total 1671	C 1044	N 281	O 339	S 7	0	0	0

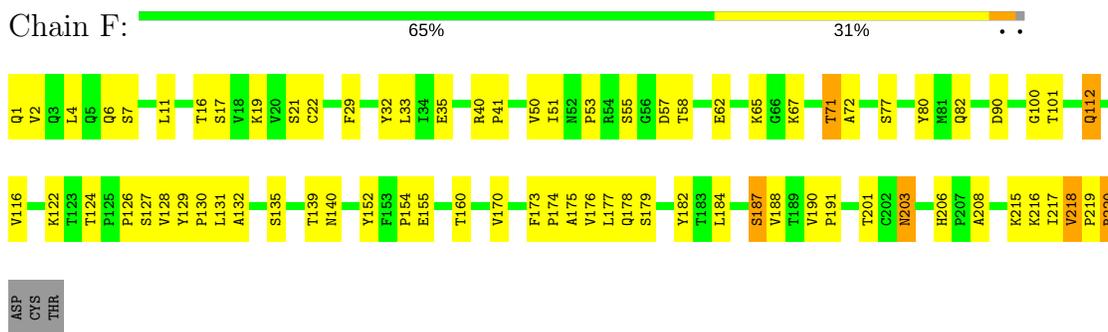
- Molecule 3 is a protein called Capsid.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	A	315	Total 2434	C 1550	N 413	O 462	S 9	0	0	0
3	B	309	Total 2397	C 1528	N 407	O 453	S 9	0	0	0

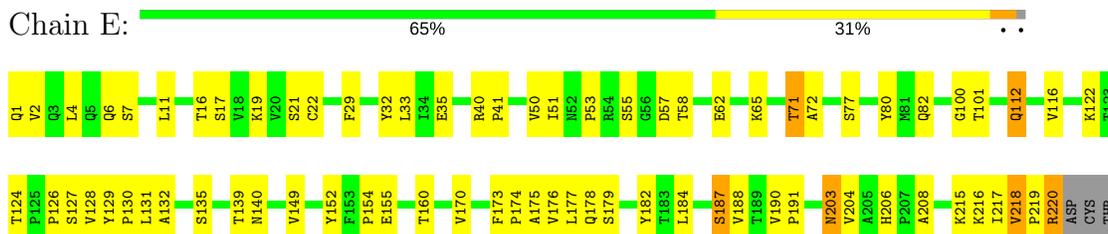
3 Residue-property plots [i](#)

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

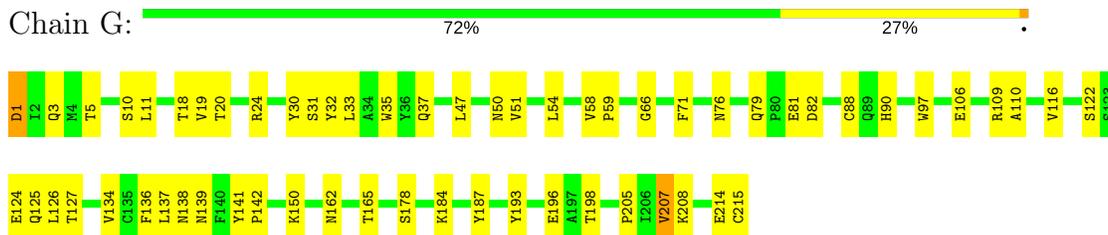
- Molecule 1: 5B18 heavy chain



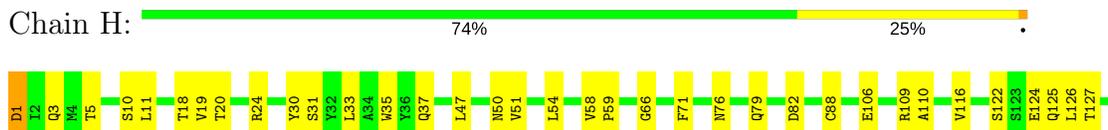
- Molecule 1: 5B18 heavy chain



- Molecule 2: 5B18 kappa chain

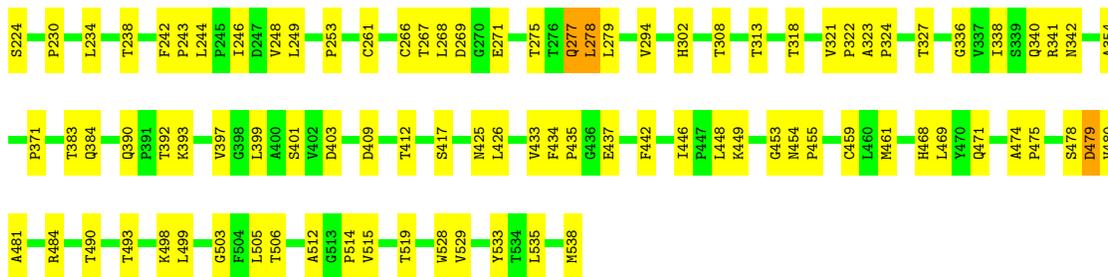


- Molecule 2: 5B18 kappa chain

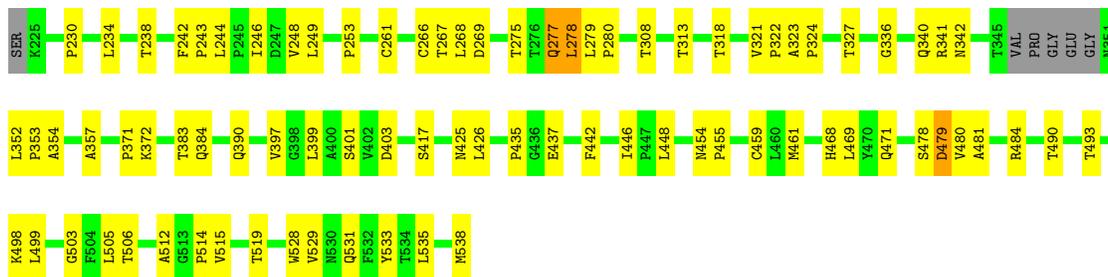




• Molecule 3: Capsid



• Molecule 3: Capsid



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 2 2	Depositor
Cell constants a, b, c, α , β , γ	145.48Å 145.48Å 216.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.54 – 3.30 31.54 – 3.30	Depositor EDS
% Data completeness (in resolution range)	86.4 (31.54-3.30) 86.6 (31.54-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.03 (at 3.31Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: dev_755)	Depositor
R, R_{free}	0.227 , 0.283 0.214 , 0.277	Depositor DCC
R_{free} test set	1532 reflections (4.97%)	DCC
Wilson B-factor (Å ²)	76.8	Xtrriage
Anisotropy	0.146	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 37.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11459	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 77.18 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 8.9713e-07. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	E	0.27	0/1683	0.48	0/2300
1	F	0.27	0/1683	0.47	0/2300
2	G	0.27	0/1712	0.45	0/2325
2	H	0.27	0/1712	0.45	0/2325
3	A	0.27	0/2505	0.47	0/3433
3	B	0.27	0/2466	0.47	0/3378
All	All	0.27	0/11761	0.47	0/16061

There are no bond length outliers.

There are no bond angle outliers.

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	E	1643	0	1615	54	0
1	F	1643	0	1615	55	0
2	G	1671	0	1592	42	0
2	H	1671	0	1592	38	1
3	A	2434	0	2356	68	1
3	B	2397	0	2322	66	0
All	All	11459	0	11092	294	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 13.

The worst 5 of 294 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:H:11:LEU:HD21	2:H:19:VAL:HG21	1.53	0.89
2:G:11:LEU:HD21	2:G:19:VAL:HG21	1.55	0.88
3:A:446:ILE:O	3:B:341:ARG:NH2	2.10	0.84
3:A:399:LEU:HG	3:A:446:ILE:HG21	1.61	0.83
1:F:1:GLN:HG3	1:F:2:VAL:H	1.43	0.82

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)
2:H:150:LYS:NZ	3:A:271:GLU:OE1[3_644]	2.11	0.09

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	E	218/223 (98%)	194 (89%)	24 (11%)	0	100	100
1	F	218/223 (98%)	194 (89%)	24 (11%)	0	100	100
2	G	213/215 (99%)	193 (91%)	20 (9%)	0	100	100
2	H	213/215 (99%)	194 (91%)	19 (9%)	0	100	100
3	A	313/315 (99%)	295 (94%)	17 (5%)	1 (0%)	44	76
3	B	305/315 (97%)	288 (94%)	16 (5%)	1 (0%)	44	76
All	All	1480/1506 (98%)	1358 (92%)	120 (8%)	2 (0%)	55	85

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	243	PRO

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Mol	Chain	Res	Type
3	B	243	PRO

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	E	187/190 (98%)	177 (95%)	10 (5%)	26 62
1	F	187/190 (98%)	178 (95%)	9 (5%)	30 65
2	G	190/190 (100%)	185 (97%)	5 (3%)	51 77
2	H	190/190 (100%)	185 (97%)	5 (3%)	51 77
3	A	270/270 (100%)	260 (96%)	10 (4%)	39 71
3	B	266/270 (98%)	257 (97%)	9 (3%)	42 73
All	All	1290/1300 (99%)	1242 (96%)	48 (4%)	39 71

5 of 48 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	203	ASN
2	H	106	GLU
3	B	471	GLN
1	E	218	VAL
2	H	1	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	471	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](#)

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](#)

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	E	220/223 (98%)	-0.41	0 100 100	56, 82, 111, 156	0
1	F	220/223 (98%)	-0.40	0 100 100	54, 82, 110, 157	0
2	G	215/215 (100%)	-0.37	0 100 100	61, 84, 116, 143	0
2	H	215/215 (100%)	-0.35	1 (0%) 90 90	60, 85, 117, 144	0
3	A	315/315 (100%)	-0.40	0 100 100	52, 77, 120, 179	0
3	B	309/315 (98%)	-0.41	0 100 100	54, 78, 116, 180	0
All	All	1494/1506 (99%)	-0.39	1 (0%) 95 96	52, 81, 116, 180	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	215	CYS	2.1

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.