

# wwPDB X-ray Structure Validation Summary Report (i)

#### Aug 12, 2023 – 01:53 PM EDT

PDB ID : 8EQA

Title : Crystal structure of human anti-N1 neuraminidase 2H08 Fab

Authors : Zhu, X.; Wilson, I.A.

Deposited on : 2022-10-07

Resolution : 2.55 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

 $\begin{array}{ccc} \text{MolProbity} & : & 4.02\text{b-}467 \\ \text{Xtriage (Phenix)} & : & 1.13 \end{array}$ 

EDS : 2.35

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac : 5.8.0158

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

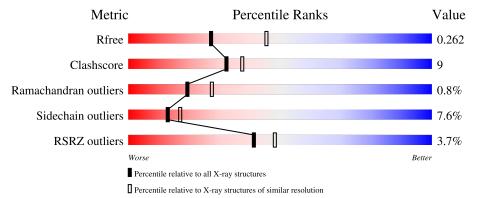
Validation Pipeline (wwPDB-VP) : 2.35

## 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\mathring{A})}) \end{array}$
$R_{free}$	130704	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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.

Mol	Chain	Length	Quality of chain					
1	A	215	78%	20%		-		
1	L	215	76%	22%		•		
2	В	225	73%	20%	•	-		
2	Н	225	75%	19%		_		



## 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6715 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 2H08 Fab light chain.

	$\mathbf{Mol}$	Chain	Residues	${f Atoms}$			ZeroOcc	AltConf	Trace		
Ī	1	т	215	Total	С	N	О	S	0	0	0
	1	ь	210	1654	1035	277	336	6	0	U	0
Ī	1	Λ	215	Total	С	N	О	S	0	0	0
	1	A	210	1654	1035	277	336	6	0	U	U

• Molecule 2 is a protein called 2H08 Fab heavy chain.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	Н	217	Total 1628	C 1035	11	O 318	S 6	0	0	0
2	R	217	Total			0	S	0	0	0
	D	211	1628	1035	269	318	6			

• Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	L	1	Total Zn 1 1	0	0
3	Н	1	Total Zn 1 1	0	0
3	A	1	Total Zn 1 1	0	0
3	В	2	$\begin{array}{cc} \text{Total} & \text{Zn} \\ 2 & 2 \end{array}$	0	0

• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	L	20	Total O 20 20	0	0

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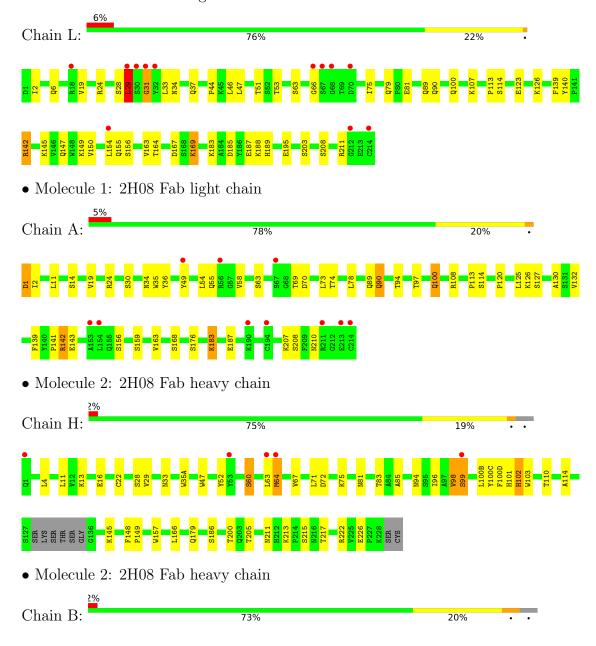
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	Н	42	Total O 42 42	0	0
4	A	23	Total O 23 23	0	0
4	В	61	Total O 61 61	0	0



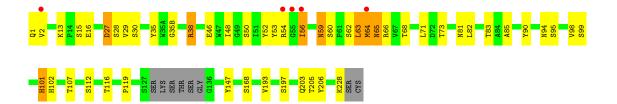
## 3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: 2H08 Fab light chain









# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants	63.56Å 110.57Å 130.68Å	Donositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.56 - 2.55	Depositor
rtesolution (A)	45.56 - 2.54	EDS
% Data completeness	97.5 (45.56-2.55)	Depositor
(in resolution range)	97.5 (45.56-2.54)	EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.18	Depositor
$< I/\sigma(I) > 1$	1.34 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
D D.	0.215 , 0.264	Depositor
$R, R_{free}$	0.212 , $0.262$	DCC
$R_{free}$ test set	1434 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.6	Xtriage
Anisotropy	0.804	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.31, 36.2	EDS
L-test for twinning <sup>2</sup>	$ < L >=0.46, < L^2>=0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	6715	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.82% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical values of <|L|>,  $<L^2>$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

## 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: ZN

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		
IVIOI	Chain	RMSZ $\# Z  > 5$		RMSZ	# Z  > 5	
1	A	0.32	0/1690	0.56	0/2295	
1	L	0.32	0/1690	0.57	0/2295	
2	В	0.35	0/1671	0.56	0/2287	
2	Н	0.35	0/1671	0.58	0/2287	
All	All	0.34	0/6722	0.57	0/9164	

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	A	1654	0	1611	27	0
1	L	1654	0	1611	33	0
2	В	1628	0	1599	31	0
2	Н	1628	0	1599	28	0
3	A	1	0	0	0	0
3	В	2	0	0	0	0
3	Н	1	0	0	0	0
3	L	1	0	0	0	0
4	A	23	0	0	2	0

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Mo	Chain	Non-H	H(model)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
4	В	61	0	0	2	0
4	Н	42	0	0	2	0
4	L	20	0	0	2	0
All	All	6715	0	6420	114	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 9.

The worst 5 of 114 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:A:36:TYR:HE1	1:A:89:GLN:HG2	1.43	0.81
2:B:29:VAL:HB	2:B:71:LEU:HD21	1.65	0.78
2:B:59:ASN:HB2	2:B:64:MET:HG2	1.70	0.74
2:B:94:ASN:HB3	2:B:101:HIS:HB3	1.71	0.73
1:L:6:GLN:H	1:L:100:GLN:NE2	1.87	0.72

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	Perce	entiles
1	A	$213/215 \ (99\%)$	207 (97%)	6 (3%)	0	100	100
1	L	213/215 (99%)	205 (96%)	6 (3%)	2 (1%)	17	24
2	В	213/225 (95%)	202 (95%)	9 (4%)	2 (1%)	17	24
2	Н	213/225 (95%)	199 (93%)	11 (5%)	3 (1%)	11	15
All	All	852/880 (97%)	813 (95%)	32 (4%)	7 (1%)	19	27

5 of 7 Ramachandran outliers are listed below:



Mol	Chain	Res	Type
2	Н	64	MET
2	Н	100(D)	PHE
1	L	29	LEU
1	L	31	GLY
2	Н	102	HIS

#### 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	$190/190\ (100\%)$	172 (90%)	18 (10%)	8 10
1	L	190/190 (100%)	178 (94%)	12 (6%)	18 23
2	В	$187/194\ (96\%)$	170 (91%)	17 (9%)	9 11
2	Н	$187/194\ (96\%)$	177 (95%)	10 (5%)	22 30
All	All	754/768~(98%)	697 (92%)	57 (8%)	13 17

5 of 57 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	GLN
2	В	116	THR
1	A	168	SER
2	В	112	SER
2	В	65	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
2	В	94	ASN
1	A	90	GLN
1	A	79	GLN
2	Н	179	GLN
1	A	89	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 monosaccharides in this entry.

#### 5.6 Ligand geometry (i)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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,  $95^{th}$  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>	$\# \mathrm{RSRZ}{>}2$	$\mathbf{OWAB}(\mathbf{\mathring{A}}^2)$	Q < 0.9
1	A	215/215 (100%)	0.40	10 (4%) 31 38	33, 51, 79, 124	0
1	L	215/215 (100%)	0.33	12 (5%) 24 29	28, 45, 78, 130	0
2	В	217/225~(96%)	0.10	5 (2%) 60 67	25, 35, 72, 152	0
2	Н	217/225 (96%)	0.10	5 (2%) 60 67	22, 33, 67, 122	0
All	All	864/880 (98%)	0.23	32 (3%) 41 48	22, 42, 75, 152	0

The worst 5 of 32 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	В	64	MET	8.3
2	Н	53	TYR	6.5
1	L	66	GLY	6.5
1	L	214	CYS	6.1
1	L	30	SER	5.3

#### 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 monosaccharides in this entry.

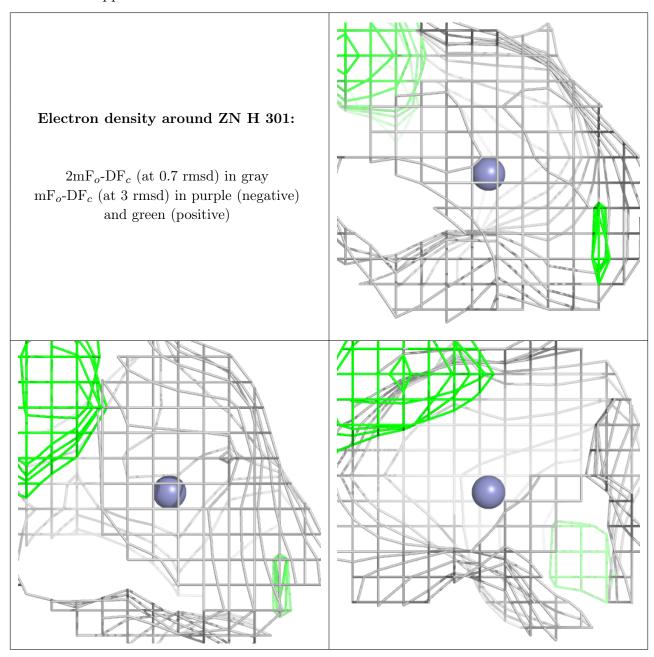
#### 6.4 Ligands (i)

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. The B-factors column lists the minimum, median,  $95^{th}$  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	$\mathbf{B} ext{-}\mathbf{factors}(\mathring{\mathbf{A}}^2)$	Q<0.9
3	ZN	Н	301	1/1	0.23	0.10	151,151,151,151	0
3	ZN	L	301	1/1	0.94	0.14	42,42,42,42	0
3	ZN	В	301	1/1	0.98	0.07	62,62,62,62	0
3	ZN	В	302	1/1	0.98	0.10	53,53,53,53	0
3	ZN	A	301	1/1	0.99	0.07	43,43,43,43	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



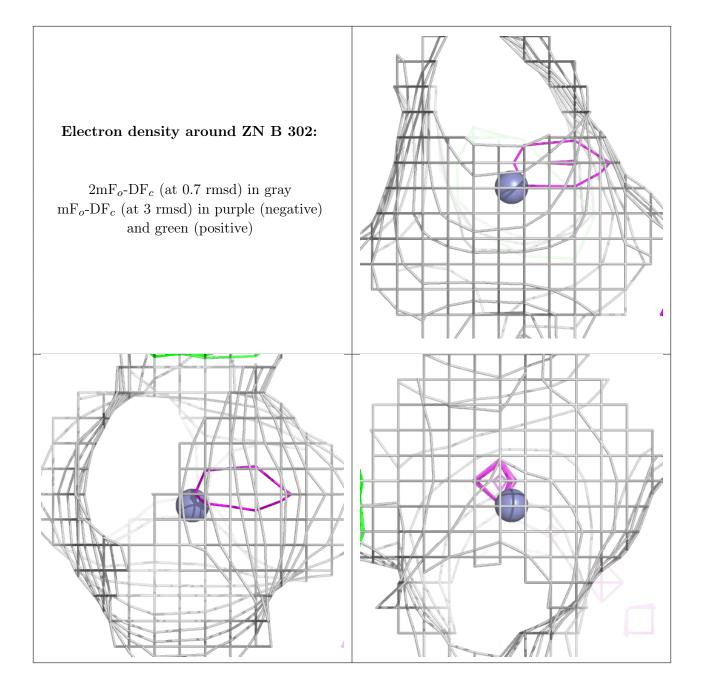


# Electron density around ZN L 301: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray $\mathrm{mF}_{o}\text{-}\mathrm{DF}_{c}$ (at 3 rmsd) in purple (negative) and green (positive)

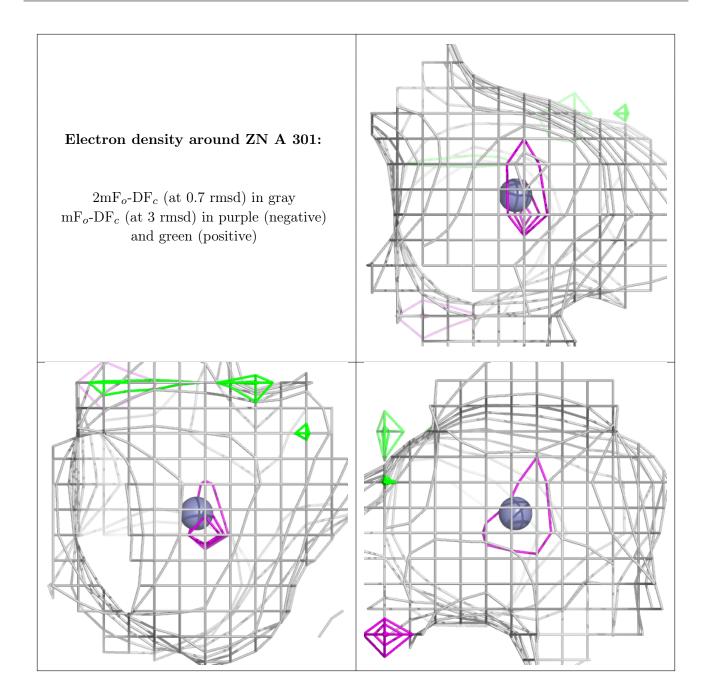


# Electron density around ZN B 301: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${ m mF}_o{ m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)









## 6.5 Other polymers (i)

There are no such residues in this entry.

