

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

#### Mar 11, 2024 – 12:36 PM JST

PDB ID	:	8IQP
Title	:	Crystal structure of Anti-PEG antibody M9 Fv-clasp fragment with PEG (co-
		crystallization with PEG3350)
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Deposited on		
Resolution	:	1.76  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:

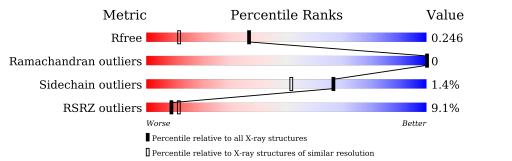
MolProbity Mogul Xtriage (Phenix) EDS	:	4.02b-467 1.8.5 (274361), CSD as541be (2020) 1.13 2.36
buster-report Percentile statistics Refmac	: : :	1.1.7 (2018) 20191225.v01 (using entries in the PDB archive December 25th 2019) 5.8.0158 7.0.044 (Gargrove)
Ideal geometry (DNA, RNA) Validation Pipeline (wwPDB-VP)		Parkinson et al. (1996) 2.36

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $X\text{-}RAY \, DIFFRACTION$ 

The reported resolution of this entry is 1.76 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
$R_{free}$	130704	$2340 \ (1.76-1.76)$
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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	L	179	<u>6%</u> 92%	• 8%
1	1	179	2% 90%	• 9%
2	Н	185	91%	• 8%
2	h	185	9%	• 8%



# 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5846 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.

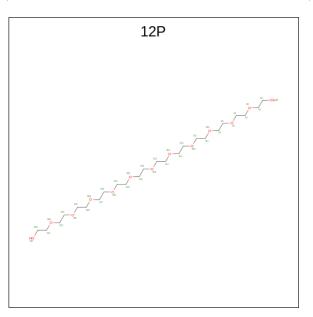
• Molecule 1 is a protein called M9 VL-SARAH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	165	Total 1316	-		0 254	S 7	0	1	0
						204	1			
1	1	163	Total	С	Ν	Ο	$\mathbf{S}$	0	1	0
1	1	105	1302	826	218	250	8	0	I	

• Molecule 2 is a protein called M9 VH-SARAH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
0	2 H 171	Total	С	Ν	0	S	0	0	0	
		1/1	1356	858	220	269	9	0	0	0
0	h	171	Total	С	Ν	0	S	0	0	0
	2 h	171	1356	858	220	269	9	U		

• Molecule 3 is DODECAETHYLENE GLYCOL (three-letter code: 12P) (formula:  $C_{24}H_{50}O_{13}$ ) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	Н	1	Total         C         O           34         22         12	0	0
3	h	1	Total         C         O           37         24         13	0	0

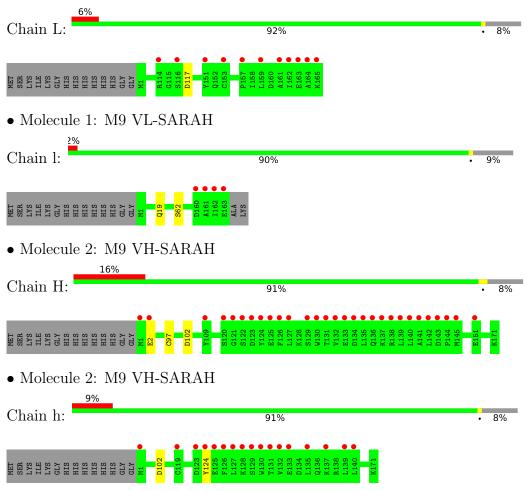
• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	L	86	Total         O           86         86	0	0
4	Н	100	Total         O           100         100	0	0
4	1	126	Total         O           126         126	0	0
4	h	133	Total O 133 133	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: M9 VL-SARAH



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	58.23Å 68.34Å 94.79Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $104.20^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	45.95 - 1.76	Depositor
Resolution (A)	45.95 - 1.76	EDS
% Data completeness	99.2 (45.95-1.76)	Depositor
(in resolution range)	99.2 (45.95 - 1.76)	EDS
R <sub>merge</sub>	(Not available)	Depositor
R <sub>sym</sub>	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.02 (at 1.76 \text{\AA})$	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
D D.	0.205 , $0.246$	Depositor
$R, R_{free}$	0.205 , $0.246$	DCC
$R_{free}$ test set	1984 reflections $(2.79\%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	33.8	Xtriage
Anisotropy	0.353	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.31 , $44.4$	EDS
L-test for twinning <sup>2</sup>	$ \langle L  \rangle = 0.51, \langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	5846	wwPDB-VP
Average B, all atoms $(Å^2)$	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The analyses of the Patterson function reveals a significant off-origin peak that is 38.94 % 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 3.4194e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.

<sup>&</sup>lt;sup>2</sup>Theoretical 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.



<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:  $12\mathrm{P}$ 

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		
	Unam	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	L	0.36	0/1347	0.58	0/1819	
1	l	0.37	0/1333	0.60	0/1799	
2	Н	0.37	0/1386	0.58	0/1872	
2	h	0.40	0/1386	0.61	0/1872	
All	All	0.37	0/5452	0.60	0/7362	

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)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

#### 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	L	164/179~(92%)	162 (99%)	2(1%)	0	100 100	

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	1	162/179~(90%)	157 (97%)	5(3%)	0	100	100
2	Н	169/185~(91%)	168 (99%)	1 (1%)	0	100	100
2	h	169/185~(91%)	168 (99%)	1 (1%)	0	100	100
All	All	664/728~(91%)	655 (99%)	9 (1%)	0	100	100

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There are no Ramachandran outliers to report.

#### 5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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	L	148/158~(94%)	147~(99%)	1 (1%)	84 75
1	1	147/158~(93%)	145~(99%)	2(1%)	67 52
2	Н	147/158~(93%)	144 (98%)	3~(2%)	55 34
2	h	147/158~(93%)	145~(99%)	2(1%)	67 52
All	All	589/632~(93%)	581 (99%)	8 (1%)	67 52

5 of 8 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	h	124	TYR
2	h	102	ASP
1	l	19	GLN
2	Н	102	ASP
1	1	62	SER

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such side chains are listed below:

Mol	Chain	Res	Type
1	1	19	GLN
1	1	152	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)

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		Bo	ond leng	$\mathbf{ths}$	В	ond ang	les
IVIOI	туре	Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
3	12P	Н	201	-	33,33,36	0.56	0	32,32,35	0.37	0
3	12P	h	201	-	36,36,36	0.55	0	35,35,35	0.43	0

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	12P	Н	201	-	-	9/31/31/34	-
3	12P	h	201	-	-	13/34/34/34	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 22 torsion outliers are listed below:

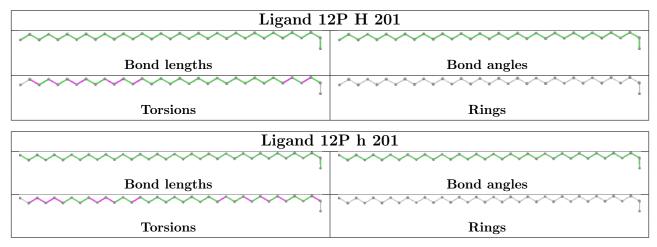


Mol	Chain	Res	Type	Atoms
3	Н	201	12P	C18-C17-O16-C15
3	Н	201	12P	C9-C8-O7-C6
3	Н	201	12P	O31-C32-C33-O34
3	h	201	12P	O4-C5-C6-O7
3	Н	201	12P	O13-C14-C15-O16

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



### 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	$\langle RSRZ \rangle$	#RSRZ>2	$OWAB(A^2)$	Q<0.9
1	L	165/179~(92%)	0.42	11 (6%) 17 23	29, 43, 74, 87	0
1	1	163/179~(91%)	0.17	4 (2%) 57 63	28, 40, 59, 80	0
2	Н	171/185~(92%)	0.68	29 (16%) 1 2	28, 41, 81, 88	0
2	h	171/185~(92%)	0.59	17 (9%) 7 10	25, 37, 80, 92	0
All	All	670/728~(92%)	0.47	61 (9%) 9 11	25, 40, 77, 92	0

The worst 5 of 61 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	h	124	TYR	11.0
2	h	131	THR	8.2
1	L	164	ALA	7.8
2	h	127	LEU	7.3
2	h	129	SER	7.2

## 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.

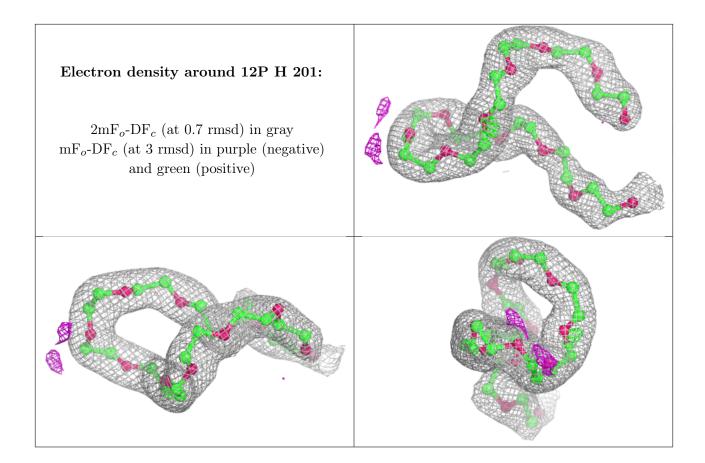


Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q < 0.9
3	12P	h	201	37/37	0.88	0.16	33,47,66,67	0
3	12P	Н	201	34/37	0.92	0.12	$34,\!43,\!55,\!59$	0

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 12P h 201: $2mF_o$ -DF <sub>c</sub> (at 0.7 rmsd) in gray $mF_o$ -DF <sub>c</sub> (at 3 rmsd) in purple (negative) and green (positive)	





### 6.5 Other polymers (i)

There are no such residues in this entry.

