

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

#### Sep 25, 2023 – 01:52 AM EDT

PDB ID : 5VPG

Title : CRYSTAL STRUCTURE OF DER P 1 COMPLEXED WITH FAB 4C1

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Deposited on : 2017-05-05

Resolution : 1.95 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 : 4.02b-467

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.35.1

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

 $Refmac \quad : \quad 5.8.0158$ 

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001)

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

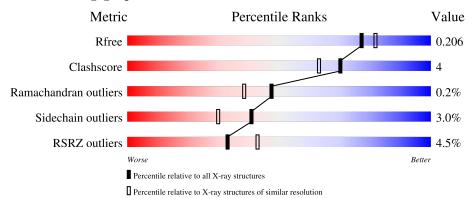
Validation Pipeline (wwPDB-VP) : 2.35.1

# 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 1.95 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\mathring{\rm A})}) \end{array}$
$R_{free}$	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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	222	91%	8% •
2	С	212	84%	14% •
3	D	255	78% 7%	15%



# 2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 5637 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 Der p 1 allergen.

I	Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
	1	A	222	Total 1788	C 1114	N 322	O 341	S 11	0	4	0

• Molecule 2 is a protein called FAB 4C1 - LIGHT CHAIN.

$\mathbf{Mol}$	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
2	С	211	Total 1674	C 1053	N 274	O 338	S 9	0	3	0

• Molecule 3 is a protein called FAB 4C1 - HEAVY CHAIN.

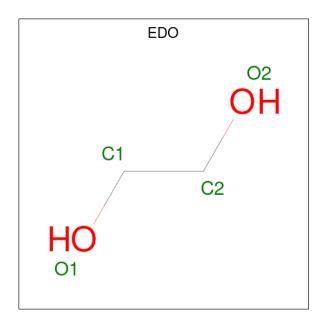
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
3	D	218	Total 1699	C 1080	N 277	O 336	S 6	0	4	0

• Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Ca 1 1	0	0

• Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).

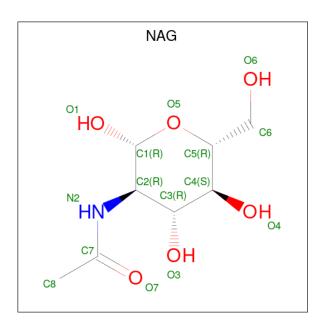




Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	С	1	Total C O 4 2 2	0	0
5	С	1	Total C O 4 2 2	0	0
5	С	1	Total C O 4 2 2	0	0

 $\bullet$  Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $\rm C_8H_{15}NO_6).$ 





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total 14	C 8	N 1	O 5	0	0

### • Molecule 7 is water.

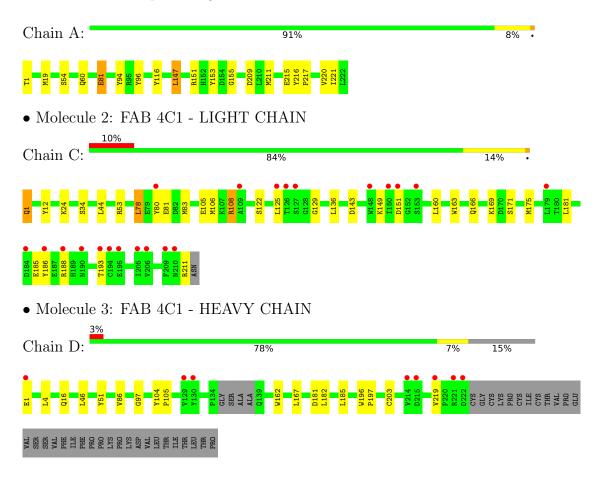
Mol	Chain	Residues	${f Atoms}$	ZeroOcc	AltConf
7	A	219	Total O 219 219	0	0
7	С	100	Total O 100 100	0	0
7	D	110	Total O 110 110	0	0



# 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: Der p 1 allergen





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	49.80Å 61.76Å 223.77Å	Donositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.50 - 1.95	Depositor
Resolution (A)	41.46 - 1.95	EDS
% Data completeness	92.7 (45.50-1.95)	Depositor
(in resolution range)	92.7 (41.46-1.95)	EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.09	Depositor
$< I/\sigma(I) > 1$	4.38 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
D D.	0.156 , 0.199	Depositor
$R, R_{free}$	0.162 , $0.206$	DCC
$R_{free}$ test set	2401 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.5	Xtriage
Anisotropy	0.139	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.34, 49.7	EDS
L-test for twinning <sup>2</sup>	$ < L > = 0.48, < L^2> = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5637	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.04% 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: CA, EDO, NAG

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	Bo	nd lengths	Bond angles		
IVIOI		RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.96	1/1831 (0.1%)	0.85	1/2490 (0.0%)	
2	С	0.78	0/1712	0.77	1/2321 (0.0%)	
3	D	0.75	0/1746	0.75	0/2397	
All	All	0.84	$1/5289 \ (0.0\%)$	0.79	2/7208 (0.0%)	

#### All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(A)	$Ideal(\AA)$
1	A	81	GLU	CG-CD	5.13	1.59	1.51

#### All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$Ideal(^{o})$
2	С	53	ARG	NE-CZ-NH2	-5.15	117.72	120.30
1	A	211	MET	CG-SD-CE	-5.02	92.17	100.20

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	1788	0	1674	10	0
2	С	1674	0	1606	21	0

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Continued	trom	mmoninonic	maaa
COHABABACA		DIEUIUU	DUIUE
0 0 1000100000			

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	1699	0	1648	7	0
4	A	1	0	0	0	0
5	A	20	0	29	1	0
5	С	12	0	18	0	0
6	A	14	0	13	0	0
7	A	219	0	0	1	0
7	С	100	0	0	4	0
7	D	110	0	0	0	0
All	All	5637	0	4988	37	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 37 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:C:83:MET:HB3	7:C:493:HOH:O	1.99	0.63
3:D:181:ASP:O	3:D:182:LEU:HD23	1.99	0.62
2:C:106:MET:CE	7:C:493:HOH:O	2.48	0.62
2:C:181:LEU:HD22	2:C:185:GLU:HG2	1.83	0.61
2:C:78:LEU:HD13	2:C:106:MET:HE1	1.86	0.58

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	ntiles
1	A	$224/222 \ (101\%)$	215 (96%)	9 (4%)	0	100	100
2	$\mathbf{C}$	$212/212\ (100\%)$	207 (98%)	4 (2%)	1 (0%)	29	17
3	D	$218/255\ (86\%)$	212 (97%)	6 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	$654/689 \; (95\%)$	634 (97%)	19 (3%)	1 (0%)	47	38

#### All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	С	151	ASP

#### 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 Outlie		Percentiles		
1	A	186/182 (102%)	183 (98%)	3 (2%)	62 58		
2	С	190/190 (100%)	180 (95%)	10 (5%)	22 10		
3	D	196/228 (86%)	191 (97%)	5 (3%)	46 36		
All	All	572/600 (95%)	554 (97%)	18 (3%)	41 28		

5 of 18 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	D	51	TYR
3	D	219	VAL
3	D	185	LEU
2	С	108	ARG
3	D	1	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

#### 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 10 ligands modelled in this entry, 1 is monoatomic - leaving 9 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	Mol Type Chain		Res Link		Во	ond leng	ths	Bond angles		
MIOI	Туре	Chain	nes	LIIIK	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	EDO	С	303	-	3,3,3	0.50	0	2,2,2	0.70	0
5	EDO	С	302	-	3,3,3	0.32	0	2,2,2	0.27	0
5	EDO	A	302	4	3,3,3	0.54	0	2,2,2	0.63	0
5	EDO	С	301	-	3,3,3	0.72	0	2,2,2	0.94	0
5	EDO	A	304	-	3,3,3	0.37	0	2,2,2	0.69	0
5	EDO	A	303	-	3,3,3	0.64	0	2,2,2	0.32	0
5	EDO	A	306	-	3,3,3	0.37	0	2,2,2	0.34	0
6	NAG	A	307	1	14,14,15	0.79	0	17,19,21	1.10	1 (5%)
5	EDO	A	305	-	3,3,3	0.32	0	2,2,2	0.38	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
5	EDO	С	303	-	-	0/1/1/1	-
5	EDO	С	302	-	-	0/1/1/1	-
5	EDO	A	302	4	-	0/1/1/1	-
5	EDO	С	301	-	-	0/1/1/1	-
5	EDO	A	304	-	-	0/1/1/1	-
5	EDO	A	303	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	A	306	-	-	1/1/1/1	-
6	NAG	A	307	1	-	0/6/23/26	0/1/1/1
5	EDO	A	305	-	-	0/1/1/1	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

$\mathbf{Mol}$	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
6	A	307	NAG	C3-C4-C5	2.10	113.99	110.24

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	306	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	306	EDO	1	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,  $95^{th}$  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 $>$	$\#\mathrm{RSRZ}{>}2$	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q<0.9
1	A	222/222 (100%)	-0.33	0 100 100	12, 19, 29, 42	1 (0%)
2	С	$211/212\ (99\%)$	0.40	21 (9%) 7 11	13, 40, 110, 133	0
3	D	218/255~(85%)	0.02	8 (3%) 41 51	15, 34, 71, 116	0
All	All	651/689 (94%)	0.02	29 (4%) 33 43	12, 25, 89, 133	1 (0%)

The worst 5 of 29 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	С	151	ASP	6.2
2	С	153	SER	5.1
2	С	125	LEU	3.9
2	С	80	TYR	3.7
3	D	221	ARG	3.4

## 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	$\operatorname{B-factors}(\mathring{\mathbf{A}}^2)$	Q<0.9
6	NAG	A	307	14/15	0.79	0.29	50,61,62,64	0
5	EDO	С	301	4/4	0.88	0.14	28,37,38,38	0
5	EDO	A	304	4/4	0.88	0.13	45,46,46,49	0
5	EDO	С	303	4/4	0.91	0.20	38,38,39,40	0
5	EDO	С	302	4/4	0.91	0.15	27,32,34,37	0
5	EDO	A	302	4/4	0.92	0.14	21,24,28,29	0
5	EDO	A	305	4/4	0.93	0.13	35,39,40,40	0
5	EDO	A	306	4/4	0.93	0.26	35,37,38,40	0
5	EDO	A	303	4/4	0.95	0.11	22,25,25,26	0
4	CA	A	301	1/1	1.00	0.05	17,17,17,17	0

# 6.5 Other polymers (i)

There are no such residues in this entry.

