



# wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 21, 2020 – 10:46 PM BST

PDB ID : 5BPG  
Title : Crystal structure of the water-soluble FraC purified starting from the trans-membrane pore  
Authors : Caaveiro, J.M.M.; Tanaka, K.; Tsumoto, K.  
Deposited on : 2015-05-28  
Resolution : 2.14 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

<https://www.wwpdb.org/validation/2017/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
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
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)	:	2.13.1



## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5867 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 Fragaceatoxin C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	177	Total	C	N	O	S	0	1	0
			1387	882	249	252	4			
1	B	177	Total	C	N	O	S	0	0	0
			1384	880	249	251	4			
1	C	177	Total	C	N	O	S	0	0	0
			1384	880	249	251	4			
1	D	177	Total	C	N	O	S	0	0	0
			1384	880	249	251	4			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	expression tag	UNP B9W5G6
B	0	MET	-	expression tag	UNP B9W5G6
C	0	MET	-	expression tag	UNP B9W5G6
D	0	MET	-	expression tag	UNP B9W5G6

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Cl	0	0
			2	2		
2	A	1	Total	Cl	0	0
			1	1		
2	D	2	Total	Cl	0	0
			2	2		
2	C	2	Total	Cl	0	0
			2	2		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	72	Total	O	0	1
			73	73		
4	B	86	Total	O	0	0
			86	86		
4	C	81	Total	O	0	0
			81	81		
4	D	63	Total	O	0	0
			63	63		

### 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: Fragaceatoxin C

Chain A:  95% . .



- Molecule 1: Fragaceatoxin C

Chain B:  92% 7% .



- Molecule 1: Fragaceatoxin C

Chain C:  96% . .



- Molecule 1: Fragaceatoxin C

Chain D:  92% 6% .



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.30Å 44.44Å 114.95Å 90.00° 92.82° 90.00°	Depositor
Resolution (Å)	47.19 – 2.14 47.19 – 2.14	Depositor EDS
% Data completeness (in resolution range)	97.1 (47.19-2.14) 97.1 (47.19-2.14)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.24 (at 2.14Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, $R_{free}$	0.210 , 0.263 0.217 , 0.267	Depositor DCC
$R_{free}$ test set	2118 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	18.4	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 14.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.145 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	5867	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.95% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: GOL, CL

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	0.65	0/1425	0.75	0/1927
1	B	0.69	0/1419	0.79	0/1919
1	C	0.62	0/1419	0.73	0/1919
1	D	0.65	0/1419	0.76	0/1919
All	All	0.65	0/5682	0.76	0/7684

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	1387	0	1358	3	0
1	B	1384	0	1353	5	0
1	C	1384	0	1353	3	0
1	D	1384	0	1353	5	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	B	6	0	8	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	6	0	8	0	0
3	D	6	0	8	0	0
4	A	73	0	0	0	0
4	B	86	0	0	0	0
4	C	81	0	0	0	0
4	D	63	0	0	1	0
All	All	5867	0	5441	16	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 1.

The worst 5 of 16 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:B:130:GLN:NE2	1:B:134:GLU:OE2	2.24	0.71
1:D:50:THR:O	4:D:301:HOH:O	2.17	0.60
1:C:141:SER:O	1:C:152:ARG:NH2	2.41	0.53
1:B:159:LYS:HD3	1:B:179:ALA:HA	1.92	0.52
1:A:37:ILE:HD13	1:A:102:VAL:HG21	1.95	0.48

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	176/180 (98%)	173 (98%)	3 (2%)	0	100	100
1	B	175/180 (97%)	173 (99%)	2 (1%)	0	100	100
1	C	175/180 (97%)	172 (98%)	3 (2%)	0	100	100
1	D	175/180 (97%)	172 (98%)	3 (2%)	0	100	100
All	All	701/720 (97%)	690 (98%)	11 (2%)	0	100	100



There are no Ramachandran outliers to report.

### 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	142/143 (99%)	142 (100%)	0	100	100
1	B	141/143 (99%)	139 (99%)	2 (1%)	67	70
1	C	141/143 (99%)	140 (99%)	1 (1%)	84	87
1	D	141/143 (99%)	138 (98%)	3 (2%)	53	54
All	All	565/572 (99%)	559 (99%)	6 (1%)	73	76

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	61	LEU
1	D	125	GLN
1	D	86	VAL
1	B	167	SER
1	D	114	SER

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

Mol	Chain	Res	Type
1	C	76	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 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 monosaccharides in this entry.

## 5.6 Ligand geometry

Of 10 ligands modelled in this entry, 7 are monoatomic - leaving 3 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	GOL	B	203	-	5,5,5	0.41	0	5,5,5	0.39	0
3	GOL	D	203	-	5,5,5	0.33	0	5,5,5	0.44	0
3	GOL	C	203	-	5,5,5	0.38	0	5,5,5	0.37	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	GOL	B	203	-	-	0/4/4/4	-
3	GOL	D	203	-	-	3/4/4/4	-
3	GOL	C	203	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	203	GOL	C1-C2-C3-O3
3	D	203	GOL	O2-C2-C3-O3
3	D	203	GOL	O1-C1-C2-O2

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<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	177/180 (98%)	-0.41	0 100 100	9, 17, 28, 36	0
1	B	177/180 (98%)	-0.35	0 100 100	10, 17, 26, 33	0
1	C	177/180 (98%)	-0.37	0 100 100	10, 18, 26, 38	0
1	D	177/180 (98%)	-0.21	0 100 100	12, 20, 31, 41	0
All	All	708/720 (98%)	-0.33	0 100 100	9, 18, 28, 41	0

There are no RSRZ outliers to report.

### 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<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	B-factors(Å <sup>2</sup> )	Q<0.9
3	GOL	D	203	6/6	0.88	0.26	32,33,34,38	0
3	GOL	B	203	6/6	0.94	0.13	17,19,20,22	0
2	CL	B	202	1/1	0.96	0.10	33,33,33,33	0
2	CL	D	202	1/1	0.97	0.04	37,37,37,37	0

*Continued on next page...*

*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	CL	D	201	1/1	0.97	0.06	38,38,38,38	0
3	GOL	C	203	6/6	0.97	0.11	20,21,21,22	0
2	CL	A	201	1/1	0.98	0.06	24,24,24,24	0
2	CL	C	201	1/1	0.98	0.05	26,26,26,26	0
2	CL	B	201	1/1	0.99	0.07	26,26,26,26	0
2	CL	C	202	1/1	1.00	0.10	26,26,26,26	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.