

wwPDB X-ray Structure Validation Summary Report (i)

Sep 2, 2023 – 05:28 PM EDT

PDB ID	:	3RGU
Title	:	Structure of Fap-NRa at pH 5.0
Authors	:	Garnett, J.A.; Matthews, S.J.
Deposited on		
Resolution	:	3.00 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:

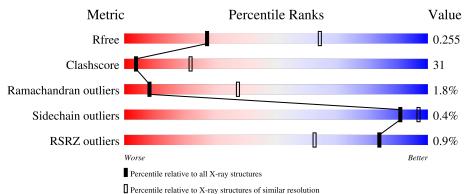
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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	А	132	.%	16% ·	33%
1	В	132	38%	28%	34%
1	С	132	.% 4 6%	17% •	36%
1	D	132	.% 41%	23% •	35%



3RGU

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 2565 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.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	88	Total	С	Ν	0	S	0	1	0
	А	00	660	407	113	137	3	0	1	
1	В	87	Total	С	Ν	0	S	0	0	0
	D	01	638	396	108	131	3	0	0	U
1	С	85	Total	С	Ν	0	S	0	0	0
	U	85	623	385	104	131	3	0	0	0
1	П	86	Total	С	Ν	0	S	0	0	0
	D	80	628	388	104	133	3	0		U

• Molecule 1 is a protein called Fimbriae-associated protein Fap1.

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	100	MET	-	expression tag	UNP Q9ZFF9
А	101	ARG	-	expression tag	UNP Q9ZFF9
А	102	GLY	-	expression tag	UNP Q9ZFF9
А	103	SER	-	expression tag	UNP Q9ZFF9
А	104	HIS	-	expression tag	UNP Q9ZFF9
А	105	HIS	-	expression tag	UNP Q9ZFF9
А	106	HIS	-	expression tag	UNP Q9ZFF9
А	107	HIS	-	expression tag	UNP Q9ZFF9
А	108	HIS	-	expression tag	UNP Q9ZFF9
А	109	HIS	-	expression tag	UNP Q9ZFF9
А	110	GLY	-	expression tag	UNP Q9ZFF9
А	111	LEU	-	expression tag	UNP Q9ZFF9
А	112	VAL	-	expression tag	UNP Q9ZFF9
А	113	PRO	-	expression tag	UNP Q9ZFF9
А	114	ARG	-	expression tag	UNP Q9ZFF9
А	115	GLY	-	expression tag	UNP Q9ZFF9
В	100	MET	-	expression tag	UNP Q9ZFF9
В	101	ARG	-	expression tag	UNP Q9ZFF9
В	102	GLY	-	expression tag	UNP Q9ZFF9
В	103	SER	-	expression tag	UNP Q9ZFF9
В	104	HIS	-	expression tag	UNP Q9ZFF9

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Chain	Residue	vious page Modelled	Actual	Comment	Reference
B	105	HIS	-	expression tag	UNP Q9ZFF9
B	106	HIS	_	expression tag	UNP Q9ZFF9
B	100	HIS	_	expression tag	UNP Q9ZFF9
B	101	HIS		expression tag	UNP Q9ZFF9
B	100	HIS		expression tag	UNP Q9ZFF9
B	105	GLY		expression tag	UNP Q9ZFF9
B	110	LEU		expression tag	UNP Q9ZFF9
B	111	VAL	-	expression tag	UNP Q9ZFF9
B	112	PRO		expression tag	UNP Q9ZFF9
B	110	ARG	-	expression tag	UNP Q9ZFF9
B	114	GLY		expression tag	UNP Q9ZFF9
C D	110	MET	-	expression tag	UNP Q9ZFF9
C	100	ARG	-	expression tag	UNP Q9ZFF9
C	101	GLY	-	expression tag	UNP Q9ZFF9
C	102	SER	-	expression tag	UNP Q9ZFF9
C	103	HIS	-		UNP Q9ZFF9
C	$\frac{104}{105}$	HIS	-	expression tag	UNP Q9ZFF9
C	105	HIS	-	expression tag	•
C			-	expression tag	UNP Q9ZFF9
C	107 108	HIS	-	expression tag	UNP Q9ZFF9
		HIS	-	expression tag	UNP Q9ZFF9
C	109	HIS	-	expression tag	UNP Q9ZFF9
C	110	GLY	-	expression tag	UNP Q9ZFF9
C	111	LEU	-	expression tag	UNP Q9ZFF9
C	112	VAL	-	expression tag	UNP Q9ZFF9
C	113	PRO	-	expression tag	UNP Q9ZFF9
C	114	ARG	-	expression tag	UNP Q9ZFF9
С	115	GLY	-	expression tag	UNP Q9ZFF9
D	100	MET	-	expression tag	UNP Q9ZFF9
D	101	ARG	-	expression tag	UNP Q9ZFF9
D	102	GLY	-	expression tag	UNP Q9ZFF9
D	103	SER	-	expression tag	UNP Q9ZFF9
D	104	HIS	-	expression tag	UNP Q9ZFF9
D	105	HIS	-	expression tag	UNP Q9ZFF9
D	106	HIS	-	expression tag	UNP Q9ZFF9
D	107	HIS	-	expression tag	UNP Q9ZFF9
D	108	HIS	-	expression tag	UNP Q9ZFF9
			1	orrespondent to a	UND OOZEEO
D	109	HIS	-	expression tag	UNP Q9ZFF9
D D	110	GLY	-	expression tag	UNP Q9ZFF9
D		GLY LEU			•
D D D D	110 111 112	GLY LEU VAL	- - - -	expression tag	UNP Q9ZFF9 UNP Q9ZFF9 UNP Q9ZFF9
D D D	110 111	GLY LEU	_	expression tag expression tag	UNP Q9ZFF9 UNP Q9ZFF9

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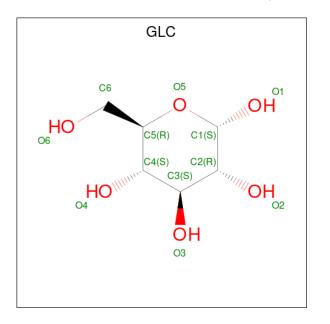
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Chain	Residue	Modelled	Actual	Comment	Reference
D	115	GLY	-	expression tag	UNP Q9ZFF9

• Molecule 2 is alpha-D-glucopyranose (three-letter code: GLC) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	С	1	Total C O 12 6 6	0	0

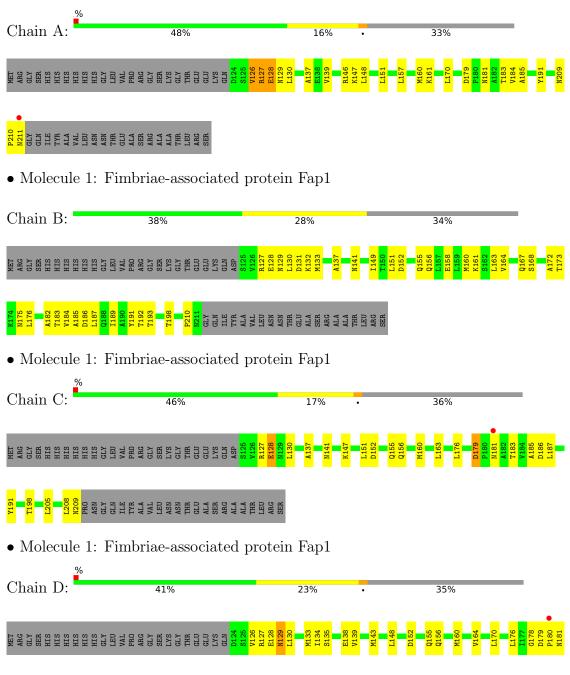
• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	2	Total O 2 2	0	0
3	В	1	Total O 1 1	0	0
3	С	1	Total O 1 1	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: Fimbriae-associated protein Fap1





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	$\begin{array}{cccccccccccccccccccccccccccccccccccc$	Depositor
Resolution (Å)	$\frac{37.43 - 3.00}{36.64 - 3.00}$	Depositor EDS
% Data completeness (in resolution range)	93.7 (37.43-3.00) 93.8 (36.64-3.00)	Depositor EDS
R _{merge} R _{sym}	0.08 (Not available)	Depositor Depositor
$< I/\sigma(I) > 1$	$2.35 (at 3.00 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.6.0104	Depositor
R, R_{free}	$\begin{array}{rrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrr$	Depositor DCC
R_{free} test set	885 reflections $(5.13%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	86.4	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.33, 81.8	EDS
L-test for $twinning^2$	$< L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	0.010 for -h,-l,-k 0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2565	wwPDB-VP
Average B, all atoms $(Å^2)$	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 16.68% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities 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: GLC

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	Bo	nd angles
	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.50	0/662	0.69	0/899
1	В	0.51	0/640	0.67	0/869
1	С	0.47	0/624	0.68	0/848
1	D	0.56	0/629	0.77	1/856~(0.1%)
All	All	0.51	0/2555	0.70	1/3472~(0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	А	0	1
1	С	0	1
All	All	0	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	D	164	VAL	CB-CA-C	-5.06	101.79	111.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	128	GLU	Peptide
1	С	128	GLU	Peptide



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	А	660	0	661	29	0
1	В	638	0	640	46	0
1	С	623	0	615	36	0
1	D	628	0	618	57	0
2	С	12	0	12	0	0
3	А	2	0	0	0	0
3	В	1	0	0	0	0
3	С	1	0	0	0	0
All	All	2565	0	2546	157	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 31.

The worst 5 of 157 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:176:LEU:HD21	1:B:187:LEU:CD1	1.66	1.25
1:C:130:LEU:HD23	1:C:191:TYR:CD1	1.71	1.24
1:C:163:LEU:HD23	1:D:160:MET:CE	1.74	1.16
1:C:130:LEU:HD21	1:C:191:TYR:HB2	1.29	1.14
1:C:130:LEU:CD2	1:C:191:TYR:HD1	1.66	1.07

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	А	87/132~(66%)	77~(88%)	8 (9%)	2(2%)	6	30
1	В	85/132~(64%)	80 (94%)	4(5%)	1 (1%)	13	48
1	С	83/132~(63%)	76~(92%)	6~(7%)	1 (1%)	13	48
1	D	84/132~(64%)	76~(90%)	6~(7%)	2(2%)	6	29
All	All	339/528~(64%)	309~(91%)	24~(7%)	6(2%)	8	37

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	128	GLU
1	А	127	ARG
1	D	184	VAL
1	А	126	VAL
1	С	179	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	Outliers	Percer	ntiles
1	А	71/110~(64%)	71~(100%)	0	100	100
1	В	67/110~(61%)	67~(100%)	0	100	100
1	С	65/110~(59%)	65 (100%)	0	100	100
1	D	66/110~(60%)	65~(98%)	1 (2%)	65	87
All	All	269/440~(61%)	268 (100%)	1 (0%)	91	97

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	129	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:



Mol	Chain	Res	Type
1	А	203	ASN
1	В	155	GLN
1	В	199	GLN
1	С	199	GLN
1	D	188	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)

1 ligand is 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	Ros	Link		ond leng			ond ang	les
WIOI	Type	Unam	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
2	GLC	С	1	-	12,12,12	0.86	0	$17,\!17,\!17$	2.05	4 (23%)

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
2	GLC	С	1	-	-	1/2/22/22	0/1/1/1



There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	С	1	GLC	C1-O5-C5	5.01	123.12	113.66
2	С	1	GLC	O5-C1-C2	4.46	118.25	110.28
2	С	1	GLC	O3-C3-C2	-2.64	104.26	110.35
2	С	1	GLC	O4-C4-C3	2.00	114.98	110.35

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	С	1	GLC	O5-C5-C6-O6

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 > 2	$\mathbf{OWAB}(\mathbf{\mathring{A}}^2)$	Q<0.9
1	А	88/132~(66%)	-0.41	1 (1%) 80 56	53, 75, 109, 141	0
1	В	87/132~(65%)	-0.25	0 100 100	54, 82, 132, 140	0
1	С	85/132~(64%)	-0.29	1 (1%) 79 54	57, 83, 140, 167	0
1	D	86/132~(65%)	-0.40	1 (1%) 79 54	47, 74, 112, 130	0
All	All	346/528~(65%)	-0.34	3 (0%) 84 63	47, 78, 130, 167	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	211	ASN	3.7
1	С	181	ASN	2.4
1	D	180	PRO	2.0

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	$B-factors(Å^2)$	Q < 0.9
2	GLC	С	1	12/12	0.92	0.46	$62,\!87,\!98,\!98$	0

6.5 Other polymers (i)

There are no such residues in this entry.

