

wwPDB X-ray Structure Validation Summary Report (i)

Oct 2, 2023 – 11:38 PM EDT

PDB ID : 6ONB

Title: Crystal Structure of the ZIG-8-RIG-5 IG1-IG1 heterodimer, monoclinic form

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Deposited on : 2019-04-20

Resolution : 1.70 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 (i)) were used in the production of this report:

MolProbity : FAILED

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

Xtriage (Phenix) : 1.13 EDS : FAILED

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

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 1.70 Å.

There are no overall percentile quality scores available for this entry.

MolProbity and EDS failed to run properly - the sequence quality summary graphics cannot be shown.



2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 4033 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 Zwei Ig domain protein zig-8.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	116	Total 930	C 579	* '	O 177	S 5	0	5	0
1	С	114	Total 892	C 556	N 163	O 168	S 5	0	2	0

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	ALA	-	expression tag	UNP G5ED00
A	20	ASP	-	expression tag	UNP G5ED00
A	21	PRO	-	expression tag	UNP G5ED00
A	138	HIS	-	expression tag	UNP G5ED00
A	139	HIS	-	expression tag	UNP G5ED00
A	140	HIS	-	expression tag	UNP G5ED00
A	141	HIS	-	expression tag	UNP G5ED00
A	142	HIS	-	expression tag	UNP G5ED00
A	143	HIS	-	expression tag	UNP G5ED00
С	19	ALA	-	expression tag	UNP G5ED00
С	20	ASP	-	expression tag	UNP G5ED00
С	21	PRO	-	expression tag	UNP G5ED00
С	138	HIS	-	expression tag	UNP G5ED00
С	139	HIS	-	expression tag	UNP G5ED00
С	140	HIS	-	expression tag	UNP G5ED00
С	141	HIS	-	expression tag	UNP G5ED00
С	142	HIS	-	expression tag	UNP G5ED00
С	143	HIS	-	expression tag	UNP G5ED00

• Molecule 2 is a protein called NeuRonal IgCAM-5.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
2	В	110	Total 896	C 561	N 162	O 169	S 4	0	4	0

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Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
2	D	109	Total 844	C 530	N 145	O 165	S 4	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
В	19	GLY	-	expression tag	UNP C6KRM7
В	20	SER	-	expression tag	UNP C6KRM7
В	131	HIS	-	expression tag	UNP C6KRM7
В	132	HIS	-	expression tag	UNP C6KRM7
В	133	HIS	-	expression tag	UNP C6KRM7
В	134	HIS	-	expression tag	UNP C6KRM7
В	135	HIS	-	expression tag	UNP C6KRM7
В	136	HIS	-	expression tag	UNP C6KRM7
D	19	GLY	-	expression tag	UNP C6KRM7
D	20	SER	-	expression tag	UNP C6KRM7
D	131	HIS	-	expression tag	UNP C6KRM7
D	132	HIS	-	expression tag	UNP C6KRM7
D	133	HIS	-	expression tag	UNP C6KRM7
D	134	HIS	-	expression tag	UNP C6KRM7
D	135	HIS	-	expression tag	UNP C6KRM7
D	136	HIS	-	expression tag	UNP C6KRM7

 \bullet Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-a cetamido-2-deoxy-beta-D-glucopyranose.

Mol	Chain	Residues	l A	\ton	ns		ZeroOcc	AltConf	Trace
3	Е	2	Total 28	C 16	N 2	O 10	0	0	0

• Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).

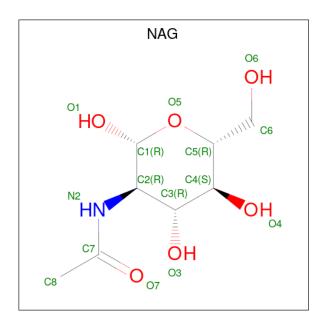




Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0
4	В	1	Total O S 5 4 1	0	0
4	В	1	Total O S 5 4 1	0	0
4	В	1	Total O S 5 4 1	0	0
4	D	1	Total O S 5 4 1	0	0

 \bullet Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $\rm C_8H_{15}NO_6).$





\mathbf{Mol}	Chain	Residues	A	tor	\mathbf{ns}		ZeroOcc	AltConf
5	D	1	Total	C 8	N 1	O 5	0	0

• Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	120	Total O 120 120	0	0
6	В	138	Total O 138 138	0	1
6	С	63	Total O 63 63	0	0
6	D	83	Total O 83 83	0	0

 $\operatorname{MolProbity}$ and EDS failed to run properly - this section is therefore empty.



3 Data and refinement statistics (i)

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	87.46Å 79.10Å 89.98Å	Depositor
a, b, c, α , β , γ	90.00° 94.22° 90.00°	Depositor
Resolution (Å)	89.73 - 1.70	Depositor
% Data completeness	96.7 (89.73-1.70)	Depositor
(in resolution range)	, , ,	•
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.14 (at 1.70Å)	Xtriage
Refinement program	PHENIX 1.15rc3_3435	Depositor
R, R_{free}	0.174 , 0.207	Depositor
Wilson B-factor (\mathring{A}^2)	27.6	Xtriage
Anisotropy	0.222	Xtriage
L-test for twinning ²	$ < L > = 0.49, < L^2> = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4033	wwPDB-VP
Average B, all atoms (\mathring{A}^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.83% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

4 Model quality (i)

4.1 Standard geometry (i)

MolProbity failed to run properly - this section is therefore empty.

4.2 Too-close contacts (i)

MolProbity failed to run properly - this section is therefore empty.

4.3 Torsion angles (i)

4.3.1 Protein backbone (i)

MolProbity failed to run properly - this section is therefore empty.

4.3.2 Protein sidechains (i)

MolProbity failed to run properly - this section is therefore empty.

4.3.3 RNA (i)

MolProbity failed to run properly - this section is therefore empty.

4.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

4.5 Carbohydrates (i)

2 monosaccharides 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 Dag Link		Link	Bo	ond leng	$ ag{ths}$	Bond angles		
MIOI	туре	Chain	Res	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
3	NAG	Е	1	3,2	14,14,15	0.47	0	17,19,21	0.48	0
3	NAG	Е	2	3	14,14,15	0.19	0	17,19,21	0.40	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	NAG	Е	1	3,2	-	2/6/23/26	0/1/1/1
3	NAG	Е	2	3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

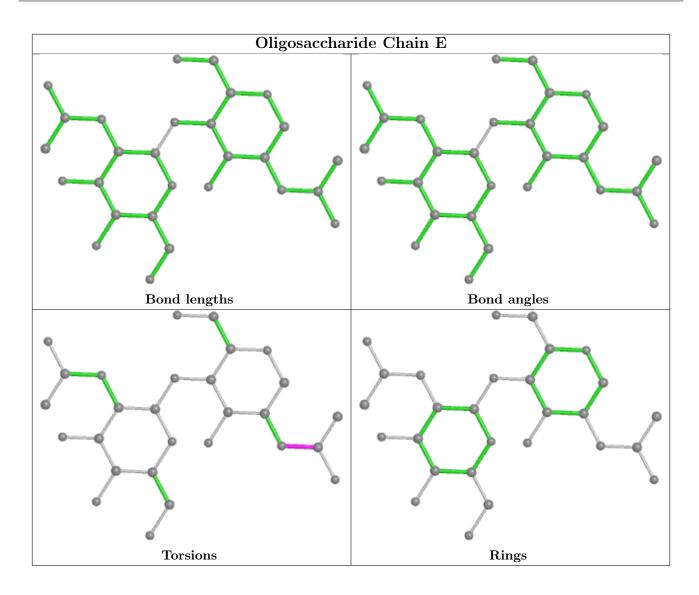
	Mol	Chain	Res	Type	Atoms
	3	Ε	1	NAG	C8-C7-N2-C2
ſ	3	Ε	1	NAG	O7-C7-N2-C2

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 oligosaccharide.





4.6 Ligand geometry (i)

6 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).

Mo	l Type	Chain	Res	Link	Bo	ond leng	$_{ m ths}$	Bond angles		
IVIO	Type				Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	В	904	-	4,4,4	0.16	0	6,6,6	0.10	0
4	SO4	В	905	-	4,4,4	0.14	0	6,6,6	0.11	0
5	NAG	D	901	2	14,14,15	0.44	0	17,19,21	0.58	0
4	SO4	В	903	-	4,4,4	0.12	0	6,6,6	0.23	0



Mol	Type Chain		Res	Link	Bo	ond leng	ths	Bond angles		
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	A	1001	-	4,4,4	0.19	0	6,6,6	0.76	0
4	SO4	D	902	-	4,4,4	0.15	0	6,6,6	0.08	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	NAG	D	901	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	D	901	NAG	C8-C7-N2-C2
5	D	901	NAG	O7-C7-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

4.7 Other polymers (i)

There are no such residues in this entry.

4.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



5 Fit of model and data (i)

5.1 Protein, DNA and RNA chains (i)

EDS failed to run properly - this section is therefore empty.

5.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS failed to run properly - this section is therefore empty.

5.3 Carbohydrates (i)

EDS failed to run properly - this section is therefore empty.

5.4 Ligands (i)

EDS failed to run properly - this section is therefore empty.

5.5 Other polymers (i)

EDS failed to run properly - this section is therefore empty.

