



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 2, 2023 – 08:48 AM EDT

PDB ID : 6MLY
Title : Bifunctional GH43-CE Bacteroides eggerthii, BACEGG_01304
Authors : Koropatkin, N.M.; Pereira, G.V.; Cann, I.
Deposited on : 2018-09-28
Resolution : 2.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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 ⓘ 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 ⓘ](#)) 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

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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 5 unique types of molecules in this entry. The entry contains 24144 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 Bifunctional GH43-CE protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	711	5555	3570	925	1032	28	0	0	0
1	B	772	6046	3871	1016	1128	31	0	0	0
1	C	774	6051	3874	1018	1128	31	0	0	0
1	D	772	6040	3869	1016	1124	31	0	0	0

There are 132 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7	GLY	-	expression tag	UNP A0A380YKU1
A	8	SER	-	expression tag	UNP A0A380YKU1
A	9	HIS	-	expression tag	UNP A0A380YKU1
A	10	MET	-	expression tag	UNP A0A380YKU1
A	11	ALA	-	expression tag	UNP A0A380YKU1
A	12	SER	-	expression tag	UNP A0A380YKU1
A	13	MET	-	expression tag	UNP A0A380YKU1
A	14	THR	-	expression tag	UNP A0A380YKU1
A	15	GLY	-	expression tag	UNP A0A380YKU1
A	16	GLY	-	expression tag	UNP A0A380YKU1
A	17	GLN	-	expression tag	UNP A0A380YKU1
A	18	GLN	-	expression tag	UNP A0A380YKU1
A	19	MET	-	expression tag	UNP A0A380YKU1
A	20	GLY	-	expression tag	UNP A0A380YKU1
A	21	ARG	-	expression tag	UNP A0A380YKU1
A	22	GLY	-	expression tag	UNP A0A380YKU1
A	23	SER	-	expression tag	UNP A0A380YKU1
A	798	TYR	-	expression tag	UNP A0A380YKU1
A	799	ARG	-	expression tag	UNP A0A380YKU1
A	800	ASP	-	expression tag	UNP A0A380YKU1
A	801	LEU	-	expression tag	UNP A0A380YKU1

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Chain	Residue	Modelled	Actual	Comment	Reference
A	802	ASN	-	expression tag	UNP A0A380YKU1
A	803	ARG	-	expression tag	UNP A0A380YKU1
A	804	ILE	-	expression tag	UNP A0A380YKU1
A	805	ARG	-	expression tag	UNP A0A380YKU1
A	806	PRO	-	expression tag	UNP A0A380YKU1
A	807	ILE	-	expression tag	UNP A0A380YKU1
A	808	LYS	-	expression tag	UNP A0A380YKU1
A	809	CYS	-	expression tag	UNP A0A380YKU1
A	810	ASP	-	expression tag	UNP A0A380YKU1
A	811	ASP	-	expression tag	UNP A0A380YKU1
A	812	ILE	-	expression tag	UNP A0A380YKU1
A	813	LYS	-	expression tag	UNP A0A380YKU1
B	7	GLY	-	expression tag	UNP A0A380YKU1
B	8	SER	-	expression tag	UNP A0A380YKU1
B	9	HIS	-	expression tag	UNP A0A380YKU1
B	10	MET	-	expression tag	UNP A0A380YKU1
B	11	ALA	-	expression tag	UNP A0A380YKU1
B	12	SER	-	expression tag	UNP A0A380YKU1
B	13	MET	-	expression tag	UNP A0A380YKU1
B	14	THR	-	expression tag	UNP A0A380YKU1
B	15	GLY	-	expression tag	UNP A0A380YKU1
B	16	GLY	-	expression tag	UNP A0A380YKU1
B	17	GLN	-	expression tag	UNP A0A380YKU1
B	18	GLN	-	expression tag	UNP A0A380YKU1
B	19	MET	-	expression tag	UNP A0A380YKU1
B	20	GLY	-	expression tag	UNP A0A380YKU1
B	21	ARG	-	expression tag	UNP A0A380YKU1
B	22	GLY	-	expression tag	UNP A0A380YKU1
B	23	SER	-	expression tag	UNP A0A380YKU1
B	798	TYR	-	expression tag	UNP A0A380YKU1
B	799	ARG	-	expression tag	UNP A0A380YKU1
B	800	ASP	-	expression tag	UNP A0A380YKU1
B	801	LEU	-	expression tag	UNP A0A380YKU1
B	802	ASN	-	expression tag	UNP A0A380YKU1
B	803	ARG	-	expression tag	UNP A0A380YKU1
B	804	ILE	-	expression tag	UNP A0A380YKU1
B	805	ARG	-	expression tag	UNP A0A380YKU1
B	806	PRO	-	expression tag	UNP A0A380YKU1
B	807	ILE	-	expression tag	UNP A0A380YKU1
B	808	LYS	-	expression tag	UNP A0A380YKU1
B	809	CYS	-	expression tag	UNP A0A380YKU1
B	810	ASP	-	expression tag	UNP A0A380YKU1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	811	ASP	-	expression tag	UNP A0A380YKU1
B	812	ILE	-	expression tag	UNP A0A380YKU1
B	813	LYS	-	expression tag	UNP A0A380YKU1
C	7	GLY	-	expression tag	UNP A0A380YKU1
C	8	SER	-	expression tag	UNP A0A380YKU1
C	9	HIS	-	expression tag	UNP A0A380YKU1
C	10	MET	-	expression tag	UNP A0A380YKU1
C	11	ALA	-	expression tag	UNP A0A380YKU1
C	12	SER	-	expression tag	UNP A0A380YKU1
C	13	MET	-	expression tag	UNP A0A380YKU1
C	14	THR	-	expression tag	UNP A0A380YKU1
C	15	GLY	-	expression tag	UNP A0A380YKU1
C	16	GLY	-	expression tag	UNP A0A380YKU1
C	17	GLN	-	expression tag	UNP A0A380YKU1
C	18	GLN	-	expression tag	UNP A0A380YKU1
C	19	MET	-	expression tag	UNP A0A380YKU1
C	20	GLY	-	expression tag	UNP A0A380YKU1
C	21	ARG	-	expression tag	UNP A0A380YKU1
C	22	GLY	-	expression tag	UNP A0A380YKU1
C	23	SER	-	expression tag	UNP A0A380YKU1
C	798	TYR	-	expression tag	UNP A0A380YKU1
C	799	ARG	-	expression tag	UNP A0A380YKU1
C	800	ASP	-	expression tag	UNP A0A380YKU1
C	801	LEU	-	expression tag	UNP A0A380YKU1
C	802	ASN	-	expression tag	UNP A0A380YKU1
C	803	ARG	-	expression tag	UNP A0A380YKU1
C	804	ILE	-	expression tag	UNP A0A380YKU1
C	805	ARG	-	expression tag	UNP A0A380YKU1
C	806	PRO	-	expression tag	UNP A0A380YKU1
C	807	ILE	-	expression tag	UNP A0A380YKU1
C	808	LYS	-	expression tag	UNP A0A380YKU1
C	809	CYS	-	expression tag	UNP A0A380YKU1
C	810	ASP	-	expression tag	UNP A0A380YKU1
C	811	ASP	-	expression tag	UNP A0A380YKU1
C	812	ILE	-	expression tag	UNP A0A380YKU1
C	813	LYS	-	expression tag	UNP A0A380YKU1
D	7	GLY	-	expression tag	UNP A0A380YKU1
D	8	SER	-	expression tag	UNP A0A380YKU1
D	9	HIS	-	expression tag	UNP A0A380YKU1
D	10	MET	-	expression tag	UNP A0A380YKU1
D	11	ALA	-	expression tag	UNP A0A380YKU1
D	12	SER	-	expression tag	UNP A0A380YKU1

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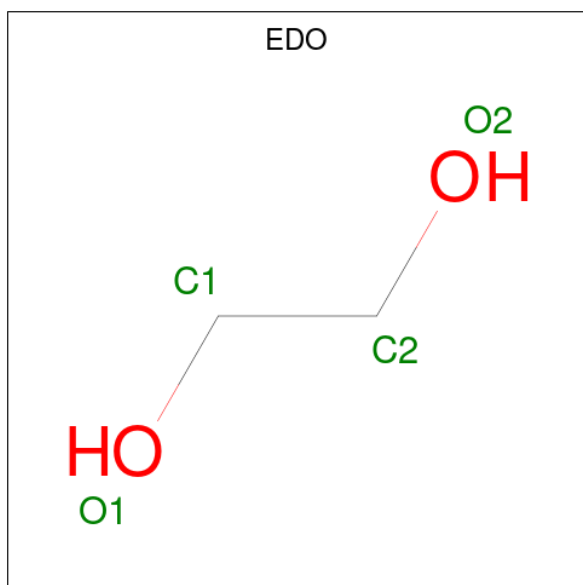
Chain	Residue	Modelled	Actual	Comment	Reference
D	13	MET	-	expression tag	UNP A0A380YKU1
D	14	THR	-	expression tag	UNP A0A380YKU1
D	15	GLY	-	expression tag	UNP A0A380YKU1
D	16	GLY	-	expression tag	UNP A0A380YKU1
D	17	GLN	-	expression tag	UNP A0A380YKU1
D	18	GLN	-	expression tag	UNP A0A380YKU1
D	19	MET	-	expression tag	UNP A0A380YKU1
D	20	GLY	-	expression tag	UNP A0A380YKU1
D	21	ARG	-	expression tag	UNP A0A380YKU1
D	22	GLY	-	expression tag	UNP A0A380YKU1
D	23	SER	-	expression tag	UNP A0A380YKU1
D	798	TYR	-	expression tag	UNP A0A380YKU1
D	799	ARG	-	expression tag	UNP A0A380YKU1
D	800	ASP	-	expression tag	UNP A0A380YKU1
D	801	LEU	-	expression tag	UNP A0A380YKU1
D	802	ASN	-	expression tag	UNP A0A380YKU1
D	803	ARG	-	expression tag	UNP A0A380YKU1
D	804	ILE	-	expression tag	UNP A0A380YKU1
D	805	ARG	-	expression tag	UNP A0A380YKU1
D	806	PRO	-	expression tag	UNP A0A380YKU1
D	807	ILE	-	expression tag	UNP A0A380YKU1
D	808	LYS	-	expression tag	UNP A0A380YKU1
D	809	CYS	-	expression tag	UNP A0A380YKU1
D	810	ASP	-	expression tag	UNP A0A380YKU1
D	811	ASP	-	expression tag	UNP A0A380YKU1
D	812	ILE	-	expression tag	UNP A0A380YKU1
D	813	LYS	-	expression tag	UNP A0A380YKU1

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	C O	0	0
			4	2 2		
2	B	1	Total	C O	0	0
			4	2 2		
2	C	1	Total	C O	0	0
			4	2 2		
2	D	1	Total	C O	0	0
			4	2 2		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	C	1	Total C O 4 2 2	0	0
3	C	1	Total C O 4 2 2	0	0
3	D	1	Total C O 4 2 2	0	0
3	D	1	Total C O 4 2 2	0	0
3	D	1	Total C O 4 2 2	0	0
3	D	1	Total C O 4 2 2	0	0

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Na 1 1	0	0
4	B	1	Total Na 1 1	0	0
4	C	1	Total Na 1 1	0	0
4	D	1	Total Na 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	58	Total O 58 58	0	0
5	B	134	Total O 134 134	0	0
5	C	92	Total O 92 92	0	0
5	D	112	Total O 112 112	0	0

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

3 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	254.73Å 93.31Å 214.04Å 90.00° 123.38° 90.00°	Depositor
Resolution (Å)	48.16 – 2.70	Depositor
% Data completeness (in resolution range)	98.8 (48.16-2.70)	Depositor
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.43 (at 2.69Å)	Xtrriage
Refinement program	REFMAC 5.8.0222	Depositor
R, R_{free}	0.220 , 0.269	Depositor
Wilson B-factor (Å ²)	44.3	Xtrriage
Anisotropy	1.150	Xtrriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	24144	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.62% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

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](#)

There are no monosaccharides in this entry.

4.6 Ligand geometry [i](#)

Of 17 ligands modelled in this entry, 4 are monoatomic - leaving 13 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	EDO	D	904	-	3,3,3	0.54	0	2,2,2	0.12	0
3	EDO	C	902	-	3,3,3	0.55	0	2,2,2	0.07	0
3	EDO	A	902	-	3,3,3	0.46	0	2,2,2	0.44	0
3	EDO	D	903	-	3,3,3	0.73	0	2,2,2	0.21	0
2	ACT	D	901	-	3,3,3	0.76	0	3,3,3	0.83	0
3	EDO	D	902	-	3,3,3	0.59	0	2,2,2	0.19	0
3	EDO	D	905	-	3,3,3	0.51	0	2,2,2	0.02	0
2	ACT	B	901	-	3,3,3	0.79	0	3,3,3	0.80	0
3	EDO	A	903	-	3,3,3	0.40	0	2,2,2	0.30	0
3	EDO	C	903	-	3,3,3	0.53	0	2,2,2	0.11	0
3	EDO	B	902	-	3,3,3	0.50	0	2,2,2	0.20	0
2	ACT	A	901	-	3,3,3	0.75	0	3,3,3	0.81	0
2	ACT	C	901	-	3,3,3	0.78	0	3,3,3	0.66	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	EDO	D	904	-	-	0/1/1/1	-
3	EDO	C	902	-	-	1/1/1/1	-
3	EDO	A	902	-	-	0/1/1/1	-
3	EDO	D	903	-	-	1/1/1/1	-
3	EDO	D	902	-	-	1/1/1/1	-
3	EDO	D	905	-	-	0/1/1/1	-
3	EDO	A	903	-	-	1/1/1/1	-
3	EDO	C	903	-	-	1/1/1/1	-
3	EDO	B	902	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	903	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	D	903	EDO	O1-C1-C2-O2
3	D	902	EDO	O1-C1-C2-O2
3	B	902	EDO	O1-C1-C2-O2
3	C	903	EDO	O1-C1-C2-O2
3	C	902	EDO	O1-C1-C2-O2

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

5.1 Protein, DNA and RNA chains

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

5.2 Non-standard residues in protein, DNA, RNA chains

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

5.3 Carbohydrates

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

5.4 Ligands

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

5.5 Other polymers

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