WORLDWIDE PDB PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report (i)

Oct 5, 2023 – 07:31 AM EDT

PDB ID	:	6VQ5
Title	:	Crystal Structure of Epiphyas postvittana Pheromone Binding Protein 3
Authors	:	Hamiaux, C.; Carraher, C.
Deposited on	:	2020-02-04
Resolution	:	2.60 Å(reported)

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 2.60 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 ⓘ

There are 3 unique types of molecules in this entry. The entry contains 1860 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 Pheromome Binding Protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	113	Total	C	N	O	S	0	0	0
			905	572	154	171	8			
1	B	117	Total	C	N	O	S	0	0	0
			930	585	158	179	8			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	124	GLU	LYS	variant	UNP A0A0K8TU48
A	143	GLU	-	expression tag	UNP A0A0K8TU48
A	144	ASN	-	expression tag	UNP A0A0K8TU48
A	145	LEU	-	expression tag	UNP A0A0K8TU48
A	146	TYR	-	expression tag	UNP A0A0K8TU48
A	147	PHE	-	expression tag	UNP A0A0K8TU48
A	148	GLN	-	expression tag	UNP A0A0K8TU48
A	149	GLY	-	expression tag	UNP A0A0K8TU48
A	150	HIS	-	expression tag	UNP A0A0K8TU48
A	151	HIS	-	expression tag	UNP A0A0K8TU48
A	152	HIS	-	expression tag	UNP A0A0K8TU48
A	153	HIS	-	expression tag	UNP A0A0K8TU48
A	154	HIS	-	expression tag	UNP A0A0K8TU48
A	155	HIS	-	expression tag	UNP A0A0K8TU48
A	156	HIS	-	expression tag	UNP A0A0K8TU48
A	157	HIS	-	expression tag	UNP A0A0K8TU48
A	158	HIS	-	expression tag	UNP A0A0K8TU48
A	159	HIS	-	expression tag	UNP A0A0K8TU48
B	124	GLU	LYS	variant	UNP A0A0K8TU48
B	143	GLU	-	expression tag	UNP A0A0K8TU48
B	144	ASN	-	expression tag	UNP A0A0K8TU48
B	145	LEU	-	expression tag	UNP A0A0K8TU48
B	146	TYR	-	expression tag	UNP A0A0K8TU48
B	147	PHE	-	expression tag	UNP A0A0K8TU48
B	148	GLN	-	expression tag	UNP A0A0K8TU48

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Chain	Residue	Modelled	Actual	Comment	Reference
B	149	GLY	-	expression tag	UNP A0A0K8TU48
B	150	HIS	-	expression tag	UNP A0A0K8TU48
B	151	HIS	-	expression tag	UNP A0A0K8TU48
B	152	HIS	-	expression tag	UNP A0A0K8TU48
B	153	HIS	-	expression tag	UNP A0A0K8TU48
B	154	HIS	-	expression tag	UNP A0A0K8TU48
B	155	HIS	-	expression tag	UNP A0A0K8TU48
B	156	HIS	-	expression tag	UNP A0A0K8TU48
B	157	HIS	-	expression tag	UNP A0A0K8TU48
B	158	HIS	-	expression tag	UNP A0A0K8TU48
B	159	HIS	-	expression tag	UNP A0A0K8TU48

- Molecule 2 is 2-(2-METHOXYETHOXY)ETHANOL (three-letter code: PG0) (formula: $C_5H_{12}O_3$).
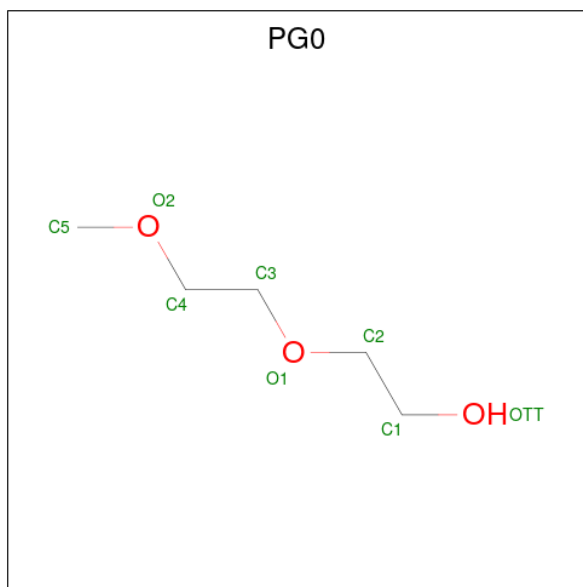
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 8	C 5	O 3	0	0
2	A	1	Total 8	C 5	O 3	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	5	Total 5	O 5	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	4	Total 4	O 4	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	P 43	Depositor
Cell constants a, b, c, α, β, γ	53.38Å 53.38Å 105.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.66 – 2.60	Depositor
% Data completeness (in resolution range)	100.0 (47.66-2.60)	Depositor
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	1.78 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.216 , 0.259	Depositor
Wilson B-factor (Å^2)	70.8	Xtriage
Anisotropy	0.425	Xtriage
L-test for twinning[2]	$<|L|> = 0.52, <L^2> = 0.36$	Xtriage
Estimated twinning fraction	0.064 for h,-k,-l	Xtriage
Total number of atoms	1860	wwPDB-VP
Average B, all atoms (Å^2)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.85% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]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.

4 Model quality ⓘ

4.1 Standard geometry ⓘ

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

4.2 Too-close contacts ⓘ

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

4.3 Torsion angles ⓘ

4.3.1 Protein backbone ⓘ

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

4.3.2 Protein sidechains ⓘ

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

4.3.3 RNA ⓘ

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

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

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

4.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

4.6 Ligand geometry ⓘ

2 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).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#$|Z| > 2$	Counts	RMSZ	#$|Z| > 2$
2	PG0	A	202	-	7,7,7	0.23	0	6,6,6	0.19	0
2	PG0	A	201	-	7,7,7	0.28	0	6,6,6	0.26	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
2	PG0	A	202	-	-	2/5/5/5	-
2	PG0	A	201	-	-	3/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	201	PG0	C1-C2-O1-C3
2	A	201	PG0	OTT-C1-C2-O1
2	A	202	PG0	C1-C2-O1-C3
2	A	202	PG0	O1-C3-C4-O2
2	A	201	PG0	O1-C3-C4-O2

There are no ring outliers.

No monomer is involved in short contacts.

4.7 Other polymers ⓘ

There are no such residues in this entry.

4.8 Polymer linkage issues ⓘ

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.