

wwPDB X-ray Structure Validation Summary Report (i)

Oct 2, 2023 – 11:24 AM EDT

PDB ID : 6MQ2

Title : De Novo Design of membrane protein-mini-eVgL membrane protein, C2221

form-2

Authors: Mravic, M.; Liu, L.; Degrado, W.F.

Deposited on : 2018-10-09

Resolution : 2.50 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 : FAILED

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

Xtriage (Phenix) : 1.13

EDS : FAILED

buster-report : 1.1.7 (2018)

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 2.50 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)

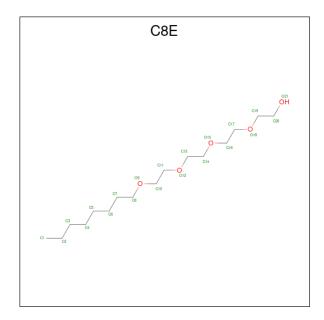
There are 3 unique types of molecules in this entry. The entry contains 1086 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 mini-eVgL membrane protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace				
1	1 A	25	Total	С	N	О	0	0	0				
1	Λ	25	207	150	29	28	0	U	0				
1	В	25	Total	С	N	О	0	0	0				
1	Ъ	25	204	149	29	26	0	0	0				
1	С	C	C	С	\mathbf{C}	25	Total	С	N	О	0	0	0
1		20	201	147	26	28	U	U	U				
1	D	27	Total	С	N	О	0	0	1				
1	D	21	212	152	31	29	0	0	1				
1	1 E	25	Total	С	N	О	0	0	0				
1		25	207	150	29	28	0		U				

• Molecule 2 is (HYDROXYETHYLOXY)TRI(ETHYLOXY)OCTANE (three-letter code: C8E) (formula: $C_{16}H_{34}O_5$).



N	lol	Chain	Residues	Atoms			ZeroOcc	AltConf
	9	Λ	1	Total	С	О	0	0
	<i>Z</i>	Α	1	12	10	2	0	0

Continued on next page...



Continued from previous page...

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 9 8 1	0	0
2	E	1	Total C 6 6	0	0
2	E	1	Total C 7 7	0	0
2	E	1	Total C O 15 12 3	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O 1 1	0	0
3	В	1	Total O 1 1	0	0
3	D	2	Total O 2 2	0	0
3	E	2	$\begin{array}{cc} \text{Total} & \text{O} \\ 2 & 2 \end{array}$	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 2 2 21	Depositor	
Cell constants	55.55Å 81.89Å 47.84Å	Depositor	
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	33.15 - 2.50	Depositor	
% Data completeness	99.0 (33.15-2.50)	Depositor	
(in resolution range)	, , ,	Dehositoi	
R_{merge}	0.16	Depositor	
R_{sym}	(Not available)	Depositor	
$< I/\sigma(I) > 1$	2.50 (at 2.51Å)	Xtriage	
Refinement program	PHENIX (1.12_2829: ???)	Depositor	
R, R_{free}	0.213 , 0.234	Depositor	
Wilson B-factor (A^2)	50.7	Xtriage	
Anisotropy	0.835	Xtriage	
L-test for twinning ²	$ < L >=0.51, < L^2>=0.35$	Xtriage	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	1086	wwPDB-VP	
Average B, all atoms (\mathring{A}^2)	66.0	wwPDB-VP	

Xtriage's analysis on translational NCS is as follows: The analyses of the Patterson function reveals a significant off-origin peak that is 24.06 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.1284e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.

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)

There are no monosaccharides in this entry.

4.6 Ligand geometry (i)

5 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	Trus	Chain	Dog Link		Res Link Bond lengths				Bond angles		
MIOI	Type	Chain	Res	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2	
2	C8E	A	102	-	8,8,20	0.16	0	7,7,19	0.48	0	
2	C8E	Е	103	-	14,14,20	0.34	0	13,13,19	0.44	0	
2	C8E	Е	101	-	5,5,20	0.27	0	4,4,19	0.40	0	
2	C8E	A	101	-	11,11,20	0.30	0	10,10,19	0.45	0	
2	C8E	Е	102	-	6,6,20	0.27	0	5,5,19	0.43	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	C8E	A	102	-	-	2/6/6/18	-
2	C8E	Е	103	-	-	8/12/12/18	-
2	C8E	Е	101	-	-	3/3/3/18	-
2	C8E	A	101	-	-	3/9/9/18	-
2	C8E	Е	102	-	-	3/4/4/18	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 19 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	Е	103	C8E	O12-C13-C14-O15
2	A	102	C8E	C3-C4-C5-C6
2	A	101	C8E	C2-C3-C4-C5
2	Е	103	C8E	C3-C4-C5-C6
2	Е	101	C8E	C2-C3-C4-C5

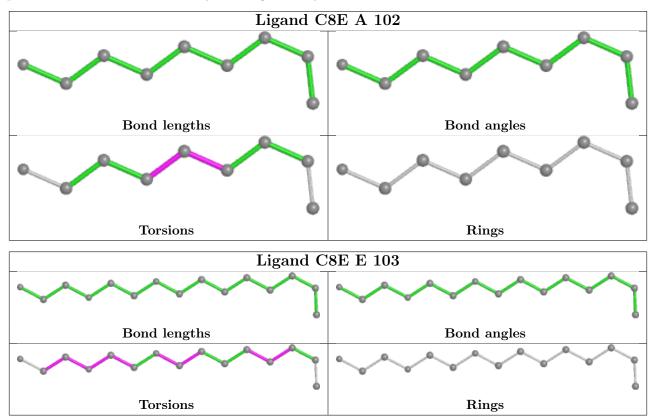
There are no ring outliers.

No monomer is involved in short contacts.

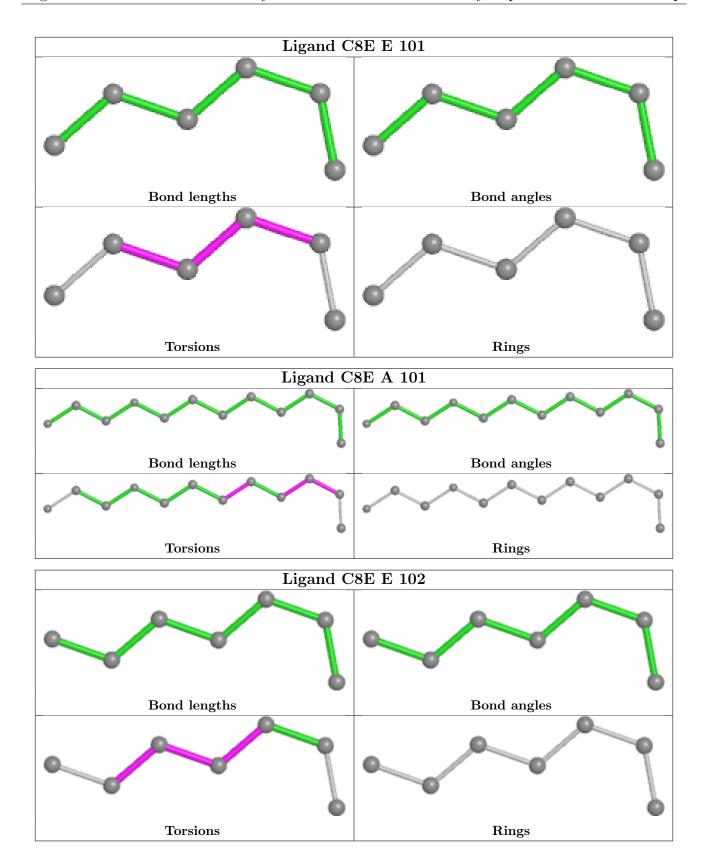
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will



also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







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.

