

# Full wwPDB X-ray Structure Validation Report (i)

Oct 2, 2023 – 02:12 AM EDT

PDB ID : 6N6I

Title : Human REXO2 bound to pGG Authors : Lormand, J.D.; Sondermann, H.

Deposited on : 2018-11-26

Resolution : 1.43 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 (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 1.43 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 7 unique types of molecules in this entry. The entry contains 3704 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.

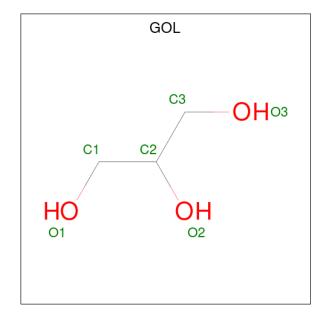
• Molecule 1 is a protein called RNA exonuclease 2 homolog, Small fragment nuclease.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	183	Total	С	N	О	S	0	19	0
1	A	100	1558	998	258	291	11	0	10	0
1	D	186	Total	С	N	О	S	0	11	0
1	Б	100	1562	998	257	296	11	0	11	0

• Molecule 2 is a RNA chain called RNA (5'-R(P\*GP\*G)-3').

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	D	9	Total C N O P	0	0					
2	D	2	47	20	10	15	2	U	U	U
9	C	9	Total	С	N	О	Р	0	0	0
2		2	47	20	10	15	2	U		U

• Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



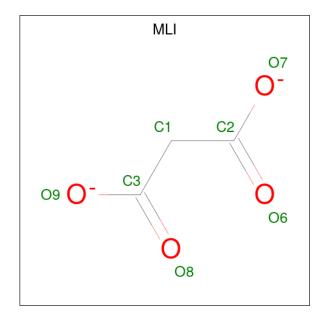


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	В	1	Total C O 6 3 3	0	0

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	D	1	Total Na 1 1	0	0
4	С	1	Total Na 1 1	0	0

• Molecule 5 is MALONATE ION (three-letter code: MLI) (formula:  $C_3H_2O_4$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	В	1	Total C O 7 3 4	0	0

 $\bullet$  Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $\mathrm{C_2H_6O_2}).$ 





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	С	1	Total C O 8 4 4	0	1

#### • Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	232	Total O 251 251	0	18
7	D	9	Total O 9 9	0	0
7	В	176	Total O 188 188	0	12
7	С	11	Total O 13 13	0	2

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	P 21 21 21	Depositor	
Cell constants	45.81Å 87.29Å 123.80Å	Depositor	
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	42.97 - 1.43	Depositor	
% Data completeness	99.0 (42.97-1.43)	Depositor	
(in resolution range)	, ,	Depositor	
$R_{merge}$	(Not available)	Depositor	
$R_{sym}$	0.10	Depositor	
$< I/\sigma(I) > 1$	1.48  (at  1.43Å)	Xtriage	
Refinement program	PHENIX 1.14_3260	Depositor	
$R, R_{free}$	0.158 , $0.178$	Depositor	
Wilson B-factor $(\mathring{A}^2)$	12.0	Xtriage	
Anisotropy	0.226	Xtriage	
L-test for twinning <sup>2</sup>	$ < L > = 0.48, < L^2> = 0.31$	Xtriage	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	3704	wwPDB-VP	
Average B, all atoms $(\mathring{A}^2)$	18.0	wwPDB-VP	

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.80% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>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.



<sup>&</sup>lt;sup>1</sup>Intensities 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)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	Trus	Chain	Res	Bond lengths			$\operatorname{gths}$	Bond angles		
MIOI	Type	Chain	nes	Link	Counts	RMSZ	# Z  > 2	Counts	RMSZ	$\mid \# Z  > 2$
5	MLI	В	302	-	6,6,6	1.38	0	7,7,7	1.23	0
6	EDO	С	701[B]	-	3,3,3	0.49	0	2,2,2	0.28	0
6	EDO	С	701[A]	-	3,3,3	0.47	0	2,2,2	0.41	0
3	GOL	В	301	-	5,5,5	0.87	0	5,5,5	1.10	0
3	GOL	A	301	-	5,5,5	0.84	0	5,5,5	0.91	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	MLI	В	302	-	-	0/4/4/4	-
6	EDO	С	701[B]	_	-	0/1/1/1	-
6	EDO	С	701[A]	-	-	1/1/1/1	-
3	GOL	В	301	-	-	0/4/4/4	-
3	GOL	A	301	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain		V -	
6	С	701[A]	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 (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.

