

Full wwPDB X-ray Structure Validation Report (i)

May 15, 2020 – 08:13 pm BST

PDB ID	:	1IRQ
Title	:	Crystal structure of omega transcriptional repressor at 1.5A resolution
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Deposited on		
Resolution	:	1.50 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 following versions of software and data (see references (1)) were used in the production of this report:

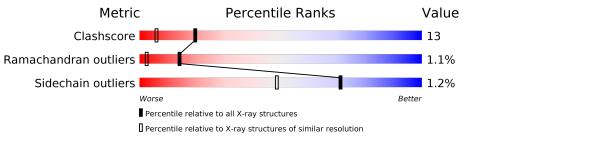
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 1.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries},{ m resolution\ range}({ m \AA}))$
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain						
1	А	71	51%	17%	32%				
1	В	71	55%	13%	• 31%				



1IRQ

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 968 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 omega transcriptional repressor.

Mol	Chain	Residues		Atc	\mathbf{ms}			ZeroOcc	AltConf	Trace
1	Λ	48	Total	С	Ν	Ο	S	0	3	0
	Л	40	412	260	74	76	2	0	J	0
1	В	49	Total	С	Ν	Ο	S	0	4	0
	D	49	431	271	76	81	3	0	4	U

• Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	67	Total O 67 67	0	0
2	В	58	Total O 58 58	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are colorcoded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

• Molecule 1: omega transcriptional repressor

Chain A:	51%	17%	32%
MET ILE VAL GLY ASN LEU GLY	GLUN GLUN LYS LYS LYS ASP ASP ASP ASP ASP ASP ASP ASP ASP AS	E43 144 N50 V51 K52 E53 E53 L59	P68 D69 L71
• Molecule	e 1: omega transcriptional repress	or	
Chain B:	55%	13% •	31%

MET VAL CLY CLY CLY CLY CLY CLY CLY ASP ASP ASP ASP ASP ASP ASP ASP ASP ASP	D23	N50	V51	K52 E53	V54	M55	T EO	R01	I63	-	P68 D69 K70	L71	
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4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 61	Depositor
Cell constants	46.10Å 46.10Å 88.30Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.60 - 1.50	Depositor
% Data completeness	97.6 (29.60-1.50)	Depositor
(in resolution range)	· · · · · ·	Depositor
R_{merge}	(Not available)	Depositor
R _{sym}	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.211 , 0.232	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	968	wwPDB-VP
Average B, all atoms $(Å^2)$	29.0	wwPDB-VP



5 Model quality (i)

5.1 Standard geometry (i)

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| > 5 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond	lengths	Bond angles			
	Ullalli	RMSZ	# Z > 5	RMSZ	# Z > 5		
1	А	0.27	0/416	0.49	0/555		
1	В	0.28	0/435	0.49	0/579		
All	All	0.27	0/851	0.49	0/1134		

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	412	0	432	12	0
1	В	431	0	446	14	0
2	А	67	0	0	2	0
2	В	58	0	0	2	0
All	All	968	0	878	22	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (22) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:ASN:HD21	1:A:52:LYS:HB3	1.53	0.74



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A 4 1	A.4.5.00	Interatomic	Clash	
Atom-1	Atom-2	distance (\AA)	overlap (Å)	
1:A:40:ILE:HG13	1:B:59[B]:LEU:HD11	1.78	0.66	
1:A:50:ASN:ND2	1:A:52:LYS:HB3	2.15	0.62	
1:A:31:ARG:O	1:B:52[A]:LYS:HG2	2.02	0.60	
1:A:50:ASN:OD1	1:A:53:GLU:HG2	2.04	0.57	
1:B:70:LYS:HB3	2:B:98:HOH:O	2.05	0.56	
1:B:50:ASN:OD1	1:B:53:GLU:HG2	2.08	0.53	
1:B:50:ASN:OD1	1:B:52[A]:LYS:HB3	2.09	0.52	
1:B:68:PRO:C	1:B:70:LYS:H	2.15	0.49	
1:B:71:LEU:HD22	1:B:71:LEU:C	2.33	0.49	
1:A:40:ILE:O	1:A:44[A]:THR:HG23	2.13	0.48	
1:B:71:LEU:N	1:B:71:LEU:HD13	2.28	0.48	
1:A:59:LEU:HD22	1:B:55[B]:MET:HE3	1.95	0.48	
1:A:59:LEU:HD22	1:B:55[B]:MET:CE	2.45	0.46	
1:A:44[B]:THR:HG23	2:A:114:HOH:O	2.15	0.46	
1:A:38:HIS:O	1:A:42:ILE:HG12	2.16	0.44	
1:A:51:VAL:HG23	2:A:95:HOH:O	2.18	0.44	
1:B:71:LEU:H	1:B:71:LEU:HD13	1.83	0.44	
1:A:68:PRO:O	1:A:69:ASP:HB2	2.19	0.43	
1:B:52[A]:LYS:HB2	2:B:100:HOH:O	2.18	0.43	
1:B:63:ILE:HG21	1:B:70:LYS:HD2	2.02	0.42	
1:B:50:ASN:OD1	1:B:52[B]:LYS:HB3	2.20	0.40	

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There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	ntiles
1	А	49/71~(69%)	47 (96%)	2(4%)	0	100	100
1	В	51/71~(72%)	47 (92%)	3~(6%)	1 (2%)	7	1
All	All	100/142~(70%)	94~(94%)	5(5%)	1 (1%)	14	3



All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	69	ASP

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	А	45/60~(75%)	45~(100%)	0	100 100	
1	В	47/60~(78%)	46~(98%)	1 (2%)	53 23	
All	All	92/120~(77%)	91~(99%)	1 (1%)	71 53	

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	В	71	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	В	47	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.



5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

