Full wwPDB X-ray Structure Validation Report (i)

Aug 20, 2023 – 09:09 PM EDT

PDB ID	:	2OL8
Title	:	The crystal structure of OspA mutant
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Deposited on	:	2007-01-18
Resolution	:	1.90 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 (1)) were used in the production of this report:

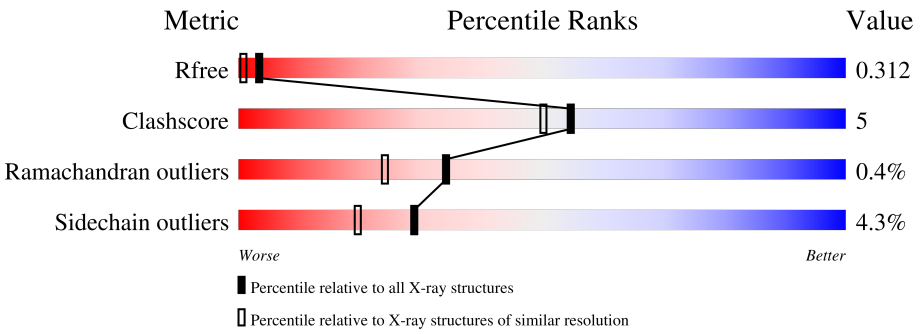
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 1.90 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	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)

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 of 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%

Mol	Chain	Length	Quality of chain
1	O	249	84% 14% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 1845 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 Outer surface protein A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	O	244	1795	1105	294	395	1	0	1	0

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
O	23	GLY	-	expression tag	UNP Q45040
O	24	SER	-	expression tag	UNP Q45040
O	25	HIS	-	expression tag	UNP Q45040
O	26	MET	-	expression tag	UNP Q45040
O	37	SER	GLU	engineered mutation	UNP Q45040
O	45	SER	GLU	engineered mutation	UNP Q45040
O	46	SER	LYS	engineered mutation	UNP Q45040
O	48	ALA	LYS	engineered mutation	UNP Q45040
O	60	ALA	LYS	engineered mutation	UNP Q45040
O	64	SER	LYS	engineered mutation	UNP Q45040
O	83	ALA	LYS	engineered mutation	UNP Q45040
O	104	SER	GLU	engineered mutation	UNP Q45040
O	107	SER	LYS	engineered mutation	UNP Q45040
O	117	ASN	LYS	engineered mutation	UNP Q45040
O	118	GLY	ASP	engineered mutation	UNP Q45040
O	?	-	LYS	deletion	UNP Q45040
O	125	ILE	PHE	engineered mutation	UNP Q45040
O	126	ILE	ASN	engineered mutation	UNP Q45040
O	127	ASP	GLU	engineered mutation	UNP Q45040
O	?	-	LYS	deletion	UNP Q45040
O	129	ILE	GLU	engineered mutation	UNP Q45040
O	130	ILE	VAL	engineered mutation	UNP Q45040
O	131	ILE	SER	engineered mutation	UNP Q45040
O	237	SER	LYS	engineered mutation	UNP Q45040
O	238	SER	GLU	engineered mutation	UNP Q45040
O	252	SER	LYS	engineered mutation	UNP Q45040

- Molecule 2 is water.

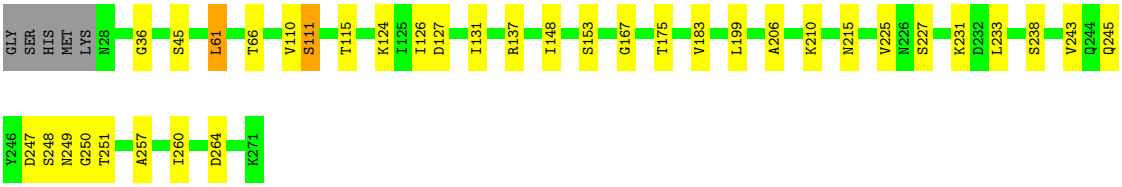
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	O	50	50	50	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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 color-coded 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.

- Molecule 1: Outer surface protein A

Chain O:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	33.36Å 53.23Å 65.19Å 90.00° 98.83° 90.00°	Depositor
Resolution (Å)	20.00 – 1.90 41.03 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (20.00-1.90) 99.5 (41.03-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.43 (at 1.81Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.232 , 0.286 0.276 , 0.312	Depositor DCC
R_{free} test set	1071 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å^2)	28.2	Xtriage
Anisotropy	0.562	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 40.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	1845	wwPDB-VP
Average B, all atoms (Å^2)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.30% of the height of the origin peak. No significant pseudotranslation is detected.*

[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.

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	
		RMSZ	$\#\|Z\|>5$	RMSZ	$\#\|Z\|>5$
1	O	0.81	0/1806	0.87	2/2435 (0.1%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	O	264	ASP	CB-CG-OD1	-7.17	111.84	118.30
1	O	137	ARG	NE-CZ-NH1	6.21	123.40	120.30

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	O	1795	0	1848	19	0
2	O	50	0	0	3	0
All	All	1845	0	1848	19	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 5.

All (19) 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:O:249:ASN:HB3	1:O:251:THR:HG23	1.57	0.87
1:O:233:LEU:HD23	1:O:243:VAL:HG22	1.70	0.74
1:O:233:LEU:CD2	1:O:243:VAL:HG22	2.22	0.68
1:O:126:ILE:HD12	1:O:131:ILE:HD13	1.84	0.59
1:O:148:ILE:HA	1:O:153:SER:O	2.12	0.49
1:O:126:ILE:O	1:O:127:ASP:HB2	2.14	0.48
1:O:183:VAL:HG13	1:O:199:LEU:HD21	1.97	0.46
1:O:215:ASN:ND2	2:O:300:HOH:O	2.50	0.43
1:O:243:VAL:CG2	1:O:260:ILE:HD11	2.48	0.43
1:O:243:VAL:HG21	1:O:260:ILE:HD11	2.00	0.43
1:O:257:ALA:O	2:O:282:HOH:O	2.21	0.43
1:O:210:LYS:HG3	1:O:225:VAL:HG22	2.00	0.43
1:O:167:GLY:HA3	1:O:175:THR:O	2.19	0.42
1:O:111:SER:OG	2:O:315:HOH:O	2.22	0.41
1:O:36:GLY:N	1:O:115:THR:HG21	2.36	0.41
1:O:231:LYS:HE3	1:O:245:GLN:OE1	2.21	0.41
1:O:249:ASN:HD22	1:O:251:THR:HG21	1.85	0.41
1:O:247:ASP:OD1	1:O:250:GLY:N	2.54	0.40
1:O:61:LEU:HD23	1:O:61:LEU:HA	1.91	0.40

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	Percentiles
1	O	243/249 (98%)	236 (97%)	6 (2%)	1 (0%)	34 24

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	O	206	ALA

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	O	210/213 (99%)	201 (96%)	9 (4%)	29 19

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	O	45	SER
1	O	61	LEU
1	O	66	THR
1	O	110	VAL
1	O	111	SER
1	O	124	LYS
1	O	227	SER
1	O	238	SER
1	O	248	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	O	249	ASN
1	O	268	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 monosaccharides 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)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.