Full wwPDB X-ray Structure Validation Report ⓘ

Nov 14, 2023 – 11:37 AM EST

PDB ID	:	8T3I
Title	:	Crystal structure of mutant exfoliative toxin C (ExhC) from Mammaliicoccus sciuri
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Deposited on	:	2023-06-07
Resolution	:	1.57 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 ⓘ 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 ⓘ) were used in the production of this report:

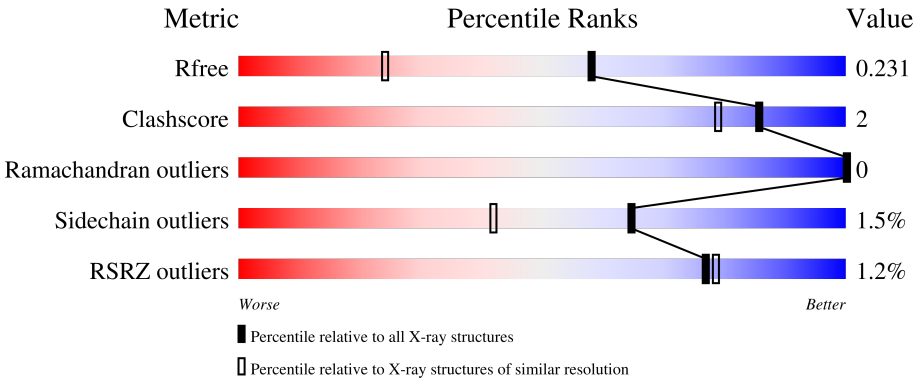
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 1.57 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	5534 (1.60-1.56)
Clashscore	141614	5861 (1.60-1.56)
Ramachandran outliers	138981	5708 (1.60-1.56)
Sidechain outliers	138945	5703 (1.60-1.56)
RSRZ outliers	127900	5431 (1.60-1.56)

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% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	258	% 86% 6% 7%
1	B	258	% 87% 6% 7%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 7625 atoms, of which 3632 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 Exfoliative toxin C.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	239	Total	C	H	N	O	S	69	0	0
			3656	1174	1810	311	360	1			
1	B	241	Total	C	H	N	O	S	69	0	0
			3686	1184	1822	313	366	1			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP F6M8N2
A	47	ALA	ARG	engineered mutation	UNP F6M8N2
A	49	PHE	ASN	engineered mutation	UNP F6M8N2
A	51	LYS	GLN	engineered mutation	UNP F6M8N2
A	89	ALA	ARG	engineered mutation	UNP F6M8N2
A	243	GLY	-	expression tag	UNP F6M8N2
A	244	GLY	-	expression tag	UNP F6M8N2
A	245	GLU	-	expression tag	UNP F6M8N2
A	246	ASN	-	expression tag	UNP F6M8N2
A	247	LEU	-	expression tag	UNP F6M8N2
A	248	TYR	-	expression tag	UNP F6M8N2
A	249	PHE	-	expression tag	UNP F6M8N2
A	250	GLN	-	expression tag	UNP F6M8N2
A	251	SER	-	expression tag	UNP F6M8N2
A	252	HIS	-	expression tag	UNP F6M8N2
A	253	HIS	-	expression tag	UNP F6M8N2
A	254	HIS	-	expression tag	UNP F6M8N2
A	255	HIS	-	expression tag	UNP F6M8N2
A	256	HIS	-	expression tag	UNP F6M8N2
A	257	HIS	-	expression tag	UNP F6M8N2
B	0	MET	-	initiating methionine	UNP F6M8N2
B	47	ALA	ARG	engineered mutation	UNP F6M8N2
B	49	PHE	ASN	engineered mutation	UNP F6M8N2
B	51	LYS	GLN	engineered mutation	UNP F6M8N2
B	89	ALA	ARG	engineered mutation	UNP F6M8N2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	243	GLY	-	expression tag	UNP F6M8N2
B	244	GLY	-	expression tag	UNP F6M8N2
B	245	GLU	-	expression tag	UNP F6M8N2
B	246	ASN	-	expression tag	UNP F6M8N2
B	247	LEU	-	expression tag	UNP F6M8N2
B	248	TYR	-	expression tag	UNP F6M8N2
B	249	PHE	-	expression tag	UNP F6M8N2
B	250	GLN	-	expression tag	UNP F6M8N2
B	251	SER	-	expression tag	UNP F6M8N2
B	252	HIS	-	expression tag	UNP F6M8N2
B	253	HIS	-	expression tag	UNP F6M8N2
B	254	HIS	-	expression tag	UNP F6M8N2
B	255	HIS	-	expression tag	UNP F6M8N2
B	256	HIS	-	expression tag	UNP F6M8N2
B	257	HIS	-	expression tag	UNP F6M8N2

- Molecule 2 is water.

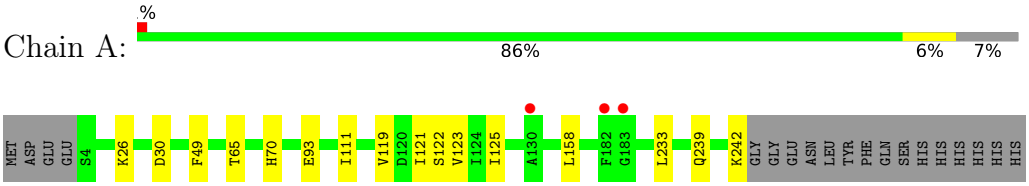
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	141	Total O 141 141	0	0
2	B	142	Total O 142 142	0	0

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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). 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.

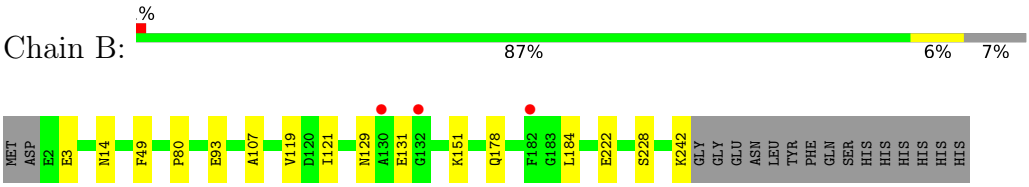
• Molecule 1: Exfoliative toxin C

Chain A:



• Molecule 1: Exfoliative toxin C

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	45.15Å 71.40Å 71.38Å 90.00° 94.70° 90.00°	Depositor
Resolution (Å)	39.55 – 1.57 39.52 – 1.57	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.55-1.57) 99.9 (39.52-1.57)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.57 (at 1.57Å)	Xtriage
Refinement program	REFMAC 5.8.0411	Depositor
R, R_{free}	0.181 , 0.223 0.192 , 0.231	Depositor DCC
R_{free} test set	3153 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	23.3	Xtriage
Anisotropy	0.454	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 37.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	7625	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.92% 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

5.1 Standard geometry

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	A	0.76	1/1888 (0.1%)	1.07	1/2551 (0.0%)
1	B	0.84	5/1906 (0.3%)	1.11	0/2575
All	All	0.80	6/3794 (0.2%)	1.09	1/5126 (0.0%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	242	LYS	CB-CG	-11.14	1.22	1.52
1	B	93	GLU	CD-OE1	8.81	1.35	1.25
1	B	3	GLU	CD-OE1	7.13	1.33	1.25
1	B	222	GLU	CD-OE2	-7.01	1.18	1.25
1	B	131	GLU	CD-OE1	6.53	1.32	1.25
1	B	222	GLU	CD-OE1	-5.28	1.19	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	30	ASP	CB-CG-OD2	-5.29	113.54	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	A	1846	1810	1803	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1864	1822	1815	5	0
2	A	141	0	0	3	0
2	B	142	0	0	1	0
All	All	3993	3632	3618	16	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 2.

All (16) 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:125:ILE:HD12	1:A:125:ILE:N	2.07	0.69
1:A:239:GLN:OE1	2:A:301:HOH:O	2.15	0.65
1:A:119:VAL:HG12	1:A:121:ILE:HG22	1.83	0.61
2:A:374:HOH:O	1:B:14:ASN:HB2	2.04	0.57
1:A:111:ILE:CD1	1:A:125:ILE:HD11	2.42	0.50
1:B:119:VAL:HG12	1:B:121:ILE:HG22	1.96	0.47
1:B:184:LEU:HG	1:B:228:SER:HB3	1.96	0.46
1:B:80:PRO:O	1:B:107:ALA:HB3	2.15	0.46
1:A:93:GLU:HG3	2:A:334:HOH:O	2.17	0.45
1:A:158:LEU:C	1:A:158:LEU:HD23	2.37	0.45
1:B:129:ASN:CB	2:B:305:HOH:O	2.65	0.45
1:A:70:HIS:HA	1:A:122:SER:OG	2.20	0.42
1:A:125:ILE:N	1:A:125:ILE:CD1	2.79	0.42
1:A:233:LEU:C	1:A:233:LEU:HD23	2.41	0.41
1:A:111:ILE:CD1	1:A:125:ILE:CD1	2.99	0.41
1:A:65:THR:HG23	1:A:123:VAL:HG13	2.04	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	A	237/258 (92%)	233 (98%)	4 (2%)	0	100	100
1	B	239/258 (93%)	233 (98%)	6 (2%)	0	100	100
All	All	476/516 (92%)	466 (98%)	10 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	A	201/218 (92%)	199 (99%)	2 (1%)	76	59
1	B	203/218 (93%)	199 (98%)	4 (2%)	55	29
All	All	404/436 (93%)	398 (98%)	6 (2%)	65	42

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	LYS
1	A	49	PHE
1	B	49	PHE
1	B	151	LYS
1	B	178	GLN
1	B	242	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	HIS
1	A	95	ASN
1	A	157	ASN
1	B	78	ASN
1	B	95	ASN
1	B	178	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	239/258 (92%)	-0.21	3 (1%)	77	78	17, 26, 46, 85	1 (0%)
1	B	241/258 (93%)	-0.22	3 (1%)	79	80	15, 26, 47, 80	1 (0%)
All	All	480/516 (93%)	-0.21	6 (1%)	77	78	15, 26, 47, 85	2 (0%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	182	PHE	8.8
1	B	130	ALA	6.1
1	B	182	PHE	6.0
1	B	132	GLY	4.2
1	A	130	ALA	3.2
1	A	183	GLY	2.1

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.