



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 26, 2024 – 07:34 AM EDT

PDB ID : 7KMM  
Title : Crystal structure of XAC1771, a novel carbohydrate acetyltransferase from *Xanthomonas citri*  
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Deposited on : 2020-11-03  
Resolution : 1.90 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
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)	:	2.37.1

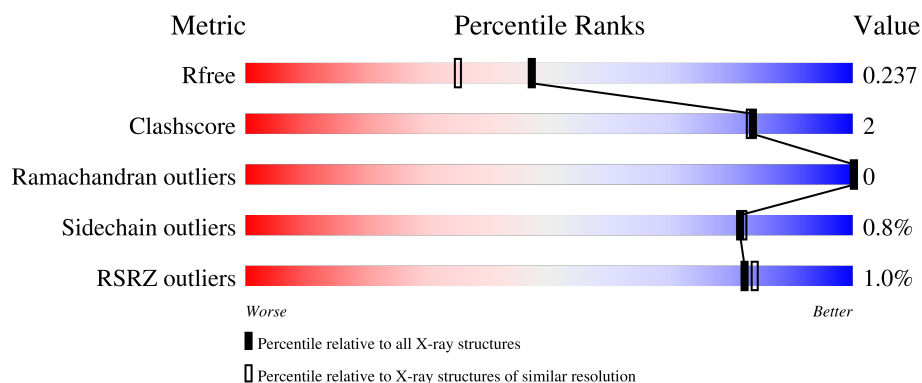
# 1 Overall quality at a glance

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)
RSRZ outliers	127900	6082 (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  $\geq 3$ , 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  $\leq 5\%$ . 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	636	<div> <div>%</div> <div> <div></div> <div>65%</div> <div>.</div> <div>32%</div> </div> </div>
1	B	636	<div> <div>%</div> <div> <div></div> <div>63%</div> <div>5%</div> <div>32%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7001 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 Sialic acid-specific 9-O-acetylerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	435	Total	C	N	O	S	0	0	0
			3314	2103	596	609	6			
1	B	432	Total	C	N	O	S	0	0	0
			3293	2090	592	605	6			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	3	MET	-	initiating methionine	UNP Q8PLM4
A	4	GLY	-	expression tag	UNP Q8PLM4
A	5	SER	-	expression tag	UNP Q8PLM4
A	6	SER	-	expression tag	UNP Q8PLM4
A	7	HIS	-	expression tag	UNP Q8PLM4
A	8	HIS	-	expression tag	UNP Q8PLM4
A	9	HIS	-	expression tag	UNP Q8PLM4
A	10	HIS	-	expression tag	UNP Q8PLM4
A	11	HIS	-	expression tag	UNP Q8PLM4
A	12	HIS	-	expression tag	UNP Q8PLM4
A	13	SER	-	expression tag	UNP Q8PLM4
A	14	SER	-	expression tag	UNP Q8PLM4
A	15	GLY	-	expression tag	UNP Q8PLM4
A	16	LEU	-	expression tag	UNP Q8PLM4
A	17	VAL	-	expression tag	UNP Q8PLM4
A	18	PRO	-	expression tag	UNP Q8PLM4
A	19	ARG	-	expression tag	UNP Q8PLM4
A	20	GLY	-	expression tag	UNP Q8PLM4
A	21	SER	-	expression tag	UNP Q8PLM4
A	22	HIS	-	expression tag	UNP Q8PLM4
A	23	MET	-	expression tag	UNP Q8PLM4
B	3	MET	-	initiating methionine	UNP Q8PLM4
B	4	GLY	-	expression tag	UNP Q8PLM4
B	5	SER	-	expression tag	UNP Q8PLM4
B	6	SER	-	expression tag	UNP Q8PLM4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	7	HIS	-	expression tag	UNP Q8PLM4
B	8	HIS	-	expression tag	UNP Q8PLM4
B	9	HIS	-	expression tag	UNP Q8PLM4
B	10	HIS	-	expression tag	UNP Q8PLM4
B	11	HIS	-	expression tag	UNP Q8PLM4
B	12	HIS	-	expression tag	UNP Q8PLM4
B	13	SER	-	expression tag	UNP Q8PLM4
B	14	SER	-	expression tag	UNP Q8PLM4
B	15	GLY	-	expression tag	UNP Q8PLM4
B	16	LEU	-	expression tag	UNP Q8PLM4
B	17	VAL	-	expression tag	UNP Q8PLM4
B	18	PRO	-	expression tag	UNP Q8PLM4
B	19	ARG	-	expression tag	UNP Q8PLM4
B	20	GLY	-	expression tag	UNP Q8PLM4
B	21	SER	-	expression tag	UNP Q8PLM4
B	22	HIS	-	expression tag	UNP Q8PLM4
B	23	MET	-	expression tag	UNP Q8PLM4

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

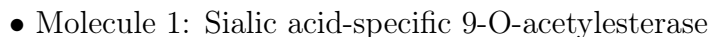
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	7	Total	Zn	0	0
			7	7		
2	B	6	Total	Zn	0	0
			6	6		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	204	Total	O	0	0
			204	204		
3	B	177	Total	O	0	0
			177	177		



- Molecule 1: Sialic acid-specific 9-O-acetylesterase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.85Å 63.50Å 88.28Å 90.00° 98.15° 90.00°	Depositor
Resolution (Å)	19.97 – 1.90 87.39 – 1.90	Depositor EDS
% Data completeness (in resolution range)	92.2 (19.97-1.90) 91.8 (87.39-1.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.83 (at 1.90Å)	Xtriage
Refinement program	PHENIX dev_3139	Depositor
R, $R_{free}$	0.190 , 0.238 0.191 , 0.237	Depositor DCC
$R_{free}$ test set	3333 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	24.3	Xtriage
Anisotropy	0.146	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 49.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7001	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.61% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, SEB

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	0.37	0/3382	0.56	0/4611
1	B	0.38	0/3361	0.55	0/4582
All	All	0.38	0/6743	0.56	0/9193

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	A	3314	0	3253	14	0
1	B	3293	0	3231	17	0
2	A	7	0	0	0	0
2	B	6	0	0	0	0
3	A	204	0	0	2	0
3	B	177	0	0	0	0
All	All	7001	0	6484	31	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.

The worst 5 of 31 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:412:HIS:HE1	3:A:949:HOH:O	1.60	0.83
1:A:412:HIS:CE1	3:A:949:HOH:O	2.41	0.65
1:B:92:GLN:HG3	1:B:97:GLU:HG2	1.82	0.62
1:A:415:GLN:HG3	1:A:453:GLU:OE1	2.07	0.53
1:B:212:LEU:HD13	1:B:403:PRO:HB3	1.91	0.53

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	430/636 (68%)	417 (97%)	13 (3%)	0	100	100
1	B	427/636 (67%)	414 (97%)	13 (3%)	0	100	100
All	All	857/1272 (67%)	831 (97%)	26 (3%)	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	328/483 (68%)	325 (99%)	3 (1%)	78	79
1	B	326/483 (68%)	324 (99%)	2 (1%)	86	87

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	654/966 (68%)	649 (99%)	5 (1%)	81	82

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	173	THR
1	A	425	TYR
1	A	484	LEU
1	B	173	THR
1	B	425	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
1	SEB	B	115	1	15,16,17	2.20	2 (13%)	15,21,23	1.58	2 (13%)
1	SEB	A	115	1	15,16,17	2.45	3 (20%)	15,21,23	1.01	1 (6%)

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
1	SEB	B	115	1	-	4/9/13/15	0/1/1/1
1	SEB	A	115	1	-	5/9/13/15	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	115	SEB	CE-SD	-7.08	1.71	1.78
1	B	115	SEB	CE-SD	-6.46	1.71	1.78
1	A	115	SEB	OG-SD	4.44	1.69	1.56
1	B	115	SEB	OG-SD	3.63	1.67	1.56
1	A	115	SEB	OD2-SD	-2.43	1.37	1.44

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	115	SEB	OD2-SD-CE	3.97	118.29	108.82
1	B	115	SEB	OG-SD-CE	-3.32	95.03	104.18
1	A	115	SEB	OG-SD-OD2	-2.35	99.00	107.39

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	115	SEB	CZ-CE-SD-OD2
1	B	115	SEB	CZ-CE-SD-OD1
1	B	115	SEB	CZ-CE-SD-OG
1	A	115	SEB	SD-CE-CZ-CH2
1	A	115	SEB	SD-CE-CZ-CH1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	115	SEB	1	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry

Of 13 ligands modelled in this entry, 13 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	434/636 (68%)	-0.12	5 (1%) 79 81	14, 24, 45, 57	0
1	B	431/636 (67%)	-0.13	4 (0%) 84 85	15, 25, 42, 66	0
All	All	865/1272 (68%)	-0.12	9 (1%) 82 84	14, 25, 43, 66	0

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	134	ALA	3.3
1	A	24(A)	VAL	3.1
1	A	604	ALA	3.0
1	A	215	ASN	3.0
1	A	397	ASP	2.7

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
1	SEB	B	115	16/17	0.88	0.18	19,40,46,51	0
1	SEB	A	115	16/17	0.92	0.16	16,39,44,46	0

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	ZN	A	706	1/1	0.66	0.11	85,85,85,85	0
2	ZN	A	701	1/1	0.93	0.04	66,66,66,66	0
2	ZN	A	705	1/1	0.95	0.09	71,71,71,71	0
2	ZN	B	703	1/1	0.95	0.07	62,62,62,62	0
2	ZN	A	704	1/1	0.97	0.04	61,61,61,61	0
2	ZN	B	702	1/1	0.98	0.07	37,37,37,37	0
2	ZN	B	705	1/1	0.98	0.05	46,46,46,46	0
2	ZN	B	701	1/1	0.99	0.04	37,37,37,37	0
2	ZN	B	704	1/1	0.99	0.03	68,68,68,68	0
2	ZN	A	702	1/1	0.99	0.06	33,33,33,33	0
2	ZN	B	706	1/1	0.99	0.04	49,49,49,49	0
2	ZN	A	703	1/1	1.00	0.03	33,33,33,33	0
2	ZN	A	707	1/1	1.00	0.11	48,48,48,48	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.